Installation Restoration Program Final First Quarter 1999-2000 Groundwater Monitoring Report

143rd Combat Communications Squadron Seattle Air National Guard Station Washington Air National Guard Seattle, Washington

January 2000



Air National Guard Andrews AFB, Maryland

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LIST OF ACRONYMS/ABBREVIATIONS

Acronym/	
Abbreviation	<u>Definition</u>
ANG	Air National Guard
ANGS	Air National Guard Station
ARAR	Applicable or relevant and appropriate requirement
ASTM	American Society for Testing and Materials
CCSQ	Combat Communications Squadron
COC	Contaminant of concern
ERM	Environmental Resources Management
IRP	Installation Restoration Program
μ g/l	Micrograms per liter
MTCA	Model Toxics Control Act
PCE	Tetrachloroethene
pН	Acidity/alkalinity
P S G	Project screening goal
QC	Quality control
RI	Remedial Investigation
TCE	Trichloroethene
VOC	Volatile organic compound
USEPA	United States Environmental Protection Agency

EXECUTIVE SUMMARY

Groundwater samples were collected in August 1999 at the Seattle Air National Guard Station in Seattle, Washington, as part of a quarterly groundwater monitoring program. Thirteen groundwater monitoring wells were sampled using low-flow purging and sampling methods. The groundwater samples were analyzed for volatile organic compounds.

The only analyte detected above project screening goals (PSGs) in August 1999 was trichloroethene (TCE). The concentration of TCE in monitoring well MW-8 near the southern Station boundary (7.2 micrograms per liter) exceeded the Washington State Model Toxics Control Act Method A Cleanup Level of 5.0 micrograms per liter. TCE was detected below the PSG in three other groundwater samples. Tetrachloroethene was detected at concentrations below the PSG in three groundwater samples. 1,1,1-trichloroethane was detected below the PSG in one groundwater sample.

Static water levels in the monitoring wells were measured prior to purging and sampling the wells. Depths to water ranged from approximately 7 to 10 feet below ground surface. The inferred direction of groundwater flow was toward the south, consistent with previous observations.

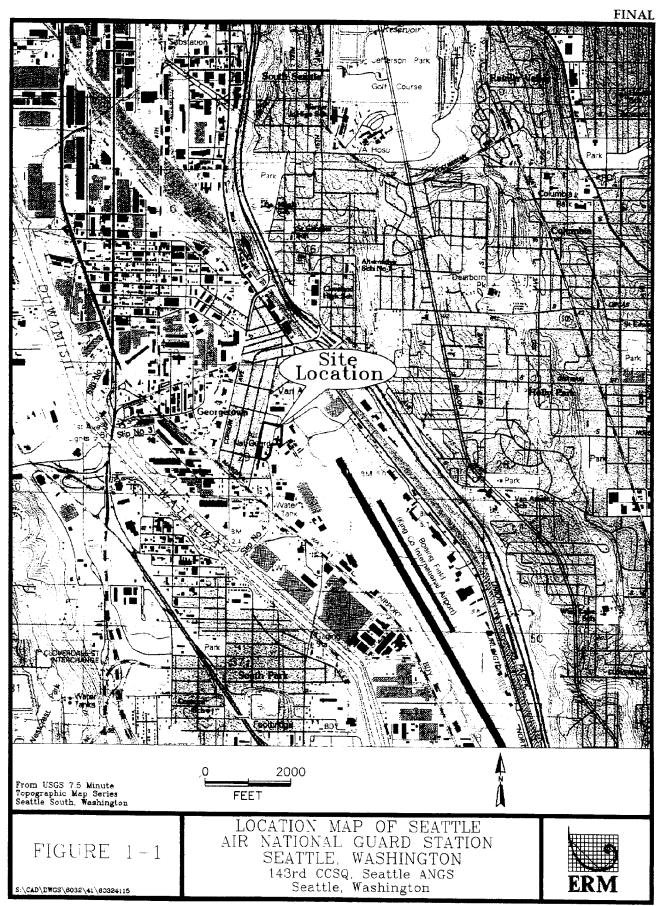
INTRODUCTION/BACKGROUND

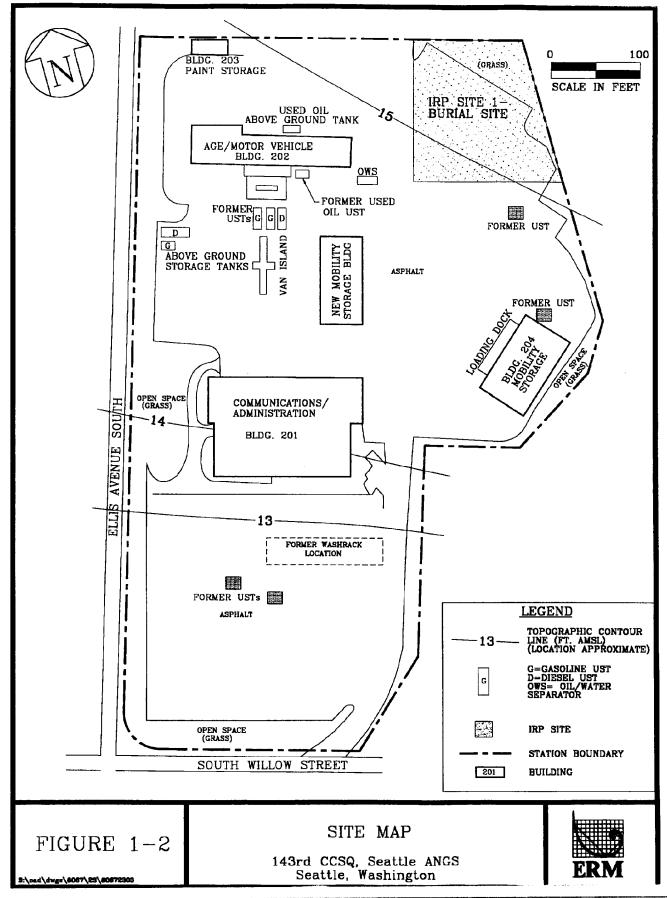
This report summarizes the methods and results of the quarterly groundwater sampling event conducted on 24-25 August 1999 at the Seattle Air National Guard Station (Seattle ANGS) in Seattle, Washington (Figure 1-1). Groundwater monitoring has been conducted at the site since September 1996 as part of the Installation Restoration Program (IRP) of the Air National Guard (ANG). Environmental Resources Management (ERM) performed the work under National Guard Bureau Contract No. DAHA90-94-D-0014, Delivery Order 0067. The Air National Guard/Installation Restoration Program Branch provided technical and project management oversight of the work.

The Seattle ANGS is at 6736 Ellis Avenue South in Seattle, Washington, and occupies approximately 7.5 acres near the north end of the King County International Airport (Boeing Field). A map of the Seattle ANGS is shown on Figure 1-2. The Seattle ANGS is currently the home of the 143rd Combat Communications Squadron (CCSQ). The mission of the 143rd CCSQ is to provide mobile communication equipment and support for airports and airfields. The facility employs approximately 129 personnel, of which 25 are full-time employees.

The Seattle ANGS was built during World War II by the War Department and was used by the United States Army Air Corps as the "Aircraft Factory School." In 1948 the property was given to King County as surplus property and was subsequently leased to the Washington ANG. On 21 April 1948, the 143rd Aircraft Control and Warning Squadron was established on the site. From May 1951 to February 1953, the 143rd was activated for recruitment purposes. During this period the unit had two C-47 aircraft. In 1960 the name of the unit was formally changed to the 143rd Communications Squadron Tributary Teams. In 1969 and 1988 the name of the unit was again changed, becoming the 143rd Mobile Communications Squadron and the 143rd CCSQ, respectively.

Currently, the Seattle ANGS property is leased from King County by the United States Air Force, which in turn licenses the property to the Washington State Military Department for ANG use.





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Three IRP investigation phases have been completed at the Seattle ANGS:

- A Preliminary Assessment, conducted by the ANG in December 1993;
- A Preliminary Assessment/Site Inspection, conducted by Operational Technologies Corporation in 1994; and
- A two-part Remedial Investigation/Feasibility Study, conducted by ERM between 1996 and 1999.

The scope and results of these IRP investigations are summarized in the Phase II Remedial Investigation (RI) and Feasibility Study Reports (ERM 1999a, 1999b). Numeric project screening goals (PSGs) were developed during the RI for use in identifying contaminants of concern (COCs) in soil and groundwater. The PSGs were derived from chemical-specific State and Federal applicable or relevant and appropriate requirements (ARARs). The derivation of PSGs and the screening process used to identify COCs are described in the Phase II RI Report (ERM 1999a).

Two chlorinated volatile organic compounds (VOCs) - trichloroethene (TCE) and tetrachloroethene (PCE) - have been detected in groundwater at concentrations above PSGs. Isolated occurrences of PCE have been detected in two background (upgradient) monitoring wells; the majority of the PCE detections have been only slightly above the Washington State Model Toxics Control Act (MTCA) Method A Groundwater Cleanup Level of 5.0 micrograms per liter (μ g/l). Dissolved TCE in groundwater is the only consistently detected COC at the Seattle ANGS that may pose a potential threat to human health or the environment. TCE has been detected at concentrations up to 83 μ g/l in shallow groundwater in the southern portion of the Station. The MTCA Method A Cleanup Level for TCE is 5.0 μ g/l.

An on-site source area for the TCE detected in groundwater has not been identified. Out of 27 soil samples analyzed for VOCs during the RI, only one was found to contain TCE. The TCE concentration reported in this sample (0.17 milligrams per kilogram) was below the MTCA Method A Soil Cleanup Level of 0.5 milligrams per kilogram. Furthermore, this soil sample was collected at the depth of the water table (approximately 10 feet below ground surface) in the southern portion of the Station, and thus may have contained TCE-impacted groundwater that biased the analytical results. Chlorinated VOCs were not detected in any of the other RI soil samples. As discussed in the Phase II RI Report (ERM 1999a), it

appears that the TCE detected in groundwater at the Seattle ANGS may be related to the groundwater contamination at the Boeing facility immediately south of the Station.

FIELD ACTIVITIES

This section provides a summary of field activities performed during the August 1999 quarterly groundwater sampling event. Groundwater sampling was conducted on 24-25 August 1999 in accordance with the procedures detailed in the 1999-2000 Groundwater Monitoring Work Plan (ERM 1999c). Figure 2-1 shows the locations of the monitoring wells.

2.1 Groundwater Sampling

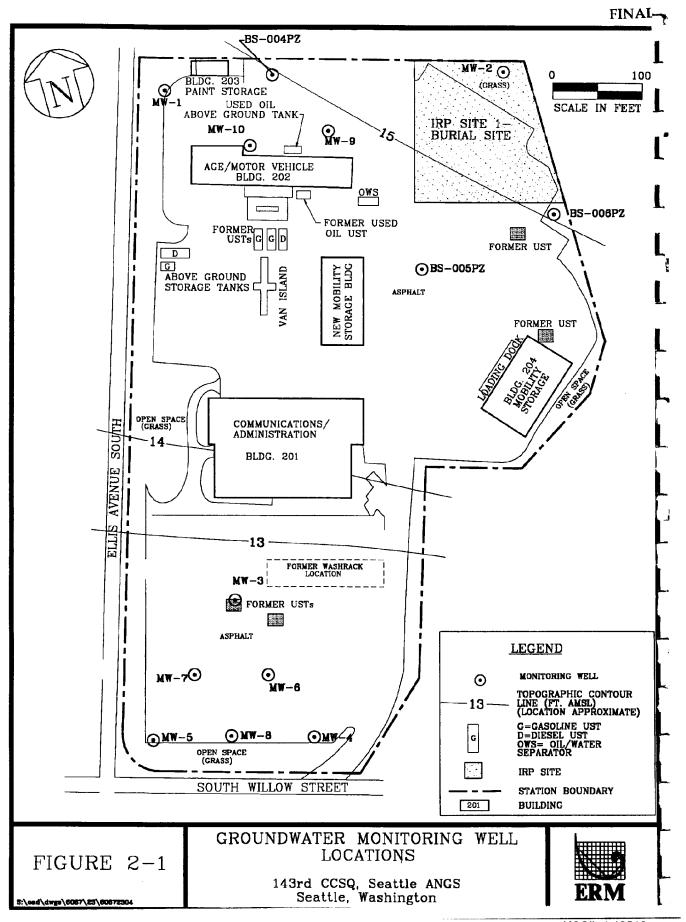
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Groundwater samples were collected and water levels measured in the following monitoring wells: BS-004PZ, BS-005PZ, BS-006PZ, and MW-1 through MW-10. Upgradient wells BS-004PZ and MW-1 (Figure 2-1) are considered background wells.

Prior to the collection of groundwater samples, static water levels in the monitoring wells were measured to within ±0.01 foot using an electronic water level indicator. Measurements were made from established reference points marked on top of each well casing. The monitoring wells were then purged and sampled using low-flow techniques. Samples were collected for analysis of VOCs.

2.1.1 Sample Collection Procedures

Using a 2-inch diameter stainless-steel submersible pump, each monitoring well was purged at a rate of less than 500 milliliters per minute. The pump was placed between the middle and top of the screened interval in each well. The temperature, acidity/alkalinity (pH), specific conductance, turbidity, and dissolved oxygen content of the purge water were monitored during well purging using an in-line flow cell and portable water quality meter. Purging continued until specific conductance and turbidity (or dissolved oxygen) stabilized to within ±10 percent, pH to within ±0.1 units, and temperature to within ±1 degree Celsius.



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After each monitoring well was purged, groundwater samples were collected using the submersible pump. The samples were collected in 40-milliliter glass vials with Teflon-lined septum lids, preserved with hydrochloric acid. Immediately following sample collection, the groundwater samples were labeled and placed in coolers containing ice. The samples were delivered to the analytical laboratory under chain of custody.

A fresh length of disposable polyethylene pump-discharge tubing was used at each monitoring well. Reusable sampling equipment was decontaminated before and after use at each well. The electronic water level indicator and the submersible pump housing were washed with an aqueous solution of Liqui-Nox (a laboratory-grade detergent) followed by a tap water rinse, a rinse with American Society for Testing and Materials (ASTM) Type II reagent-grade water, and a final spray rinse with isopropanol. The submersible pump internals were decontaminated by pumping a Liqui-Nox solution, followed by tap water and ASTM Type II water, through the pump. The ASTM Type II water was produced on site from tap water using a portable water filtration system.

2.1.2 Field QA/QC Sampling

In addition to the primary groundwater samples collected from 13 groundwater monitoring wells, the following quality assurance/quality control (QA/QC) samples were collected and/or prepared:

- One equipment rinsate blank and two field blanks. ASTM Type II water was used to prepare the rinsate blank; field blanks were prepared from tap water and ASTM Type II water. The sample identifiers for the blank samples consisted of the identifier for the primary sample collected immediately prior to the blank, followed by an "R" for the rinsate blank, an "FT" for the tap water field blank, and an "FA" for the ASTM Type II water field blank.
- One duplicate sample, from monitoring well MW-8. The sample identifier for the field duplicate sample was the same as the associated primary sample, followed by a "D".
- One sample for matrix spike/matrix spike duplicate analysis. The sample collected from monitoring well MW-8 was designated for the matrix spike/matrix spike duplicate analysis.

 One trip blank. The trip blank was prepared by the laboratory using reagent water, and accompanied the sample containers in the cooler in transit to the field and back to the laboratory.

2.1.3 Groundwater Sample Analyses

The groundwater samples were analyzed for VOCs using United States Environmental Protection Agency (USEPA) Method 8260. MultiChem Analytical Services in Renton, Washington performed the analyses. A summary of the samples submitted for laboratory analysis is shown on Table 2-1.

2.2 Investigation-Derived Waste Management

Purge water and decontamination water was collected and stored in 55-gallon drums. The groundwater analytical results were used to designate the drum contents as dangerous or non-dangerous waste in accordance with Washington State Dangerous Waste Regulations, Washington Administrative Code Chapter 173-303.

The purge water and decontamination water was determined not to be a Washington-defined dangerous waste. Accordingly, the purge water and decontamination water can be discharged to the sanitary sewer if approved by the local publicly owned treatment works (King County Department of Natural Resources, Wastewater Division). Alternatively, the purge water and decontamination water may be disposed at a facility that is permitted to receive and dispose of industrial wastewater, or at a permitted dangerous waste treatment, storage, and disposal facility.

Matrix	Sampling Method	Fiéid Parameters	Laboratory Parameters	Analytical Method	Primary Sample Analyses	* Trip * Blank	Q/ Rinsate Blank	ZQC Sar Pield Blank		MS/MSD	Total Laboratory Analyses
Groundwater	Low-flow sampling; 13 MWs, quarterly for 1 year	S.C., Turbidity, pH, Temperature, D.O.	VOCs	USEPA 8260	13	1	1	2	1	1	19

Notes:

VOC = Volatile organic compound
QA/QC = Quality assurance/quality control
DUP = Duplicate sample
pH = Acidity/alkalinity

S.C. = Specific conductance
D.O. = Dissolved oxygen content
USEPA = United States Environmental Protection Agency
MS/MSD = Matrix spike/matrix spike duplicate
MW = Monitoring well

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RESULTS

This section summarizes the results of the August 1999 groundwater sampling event at the Seattle ANGS.

3.1 Groundwater Level Data Results

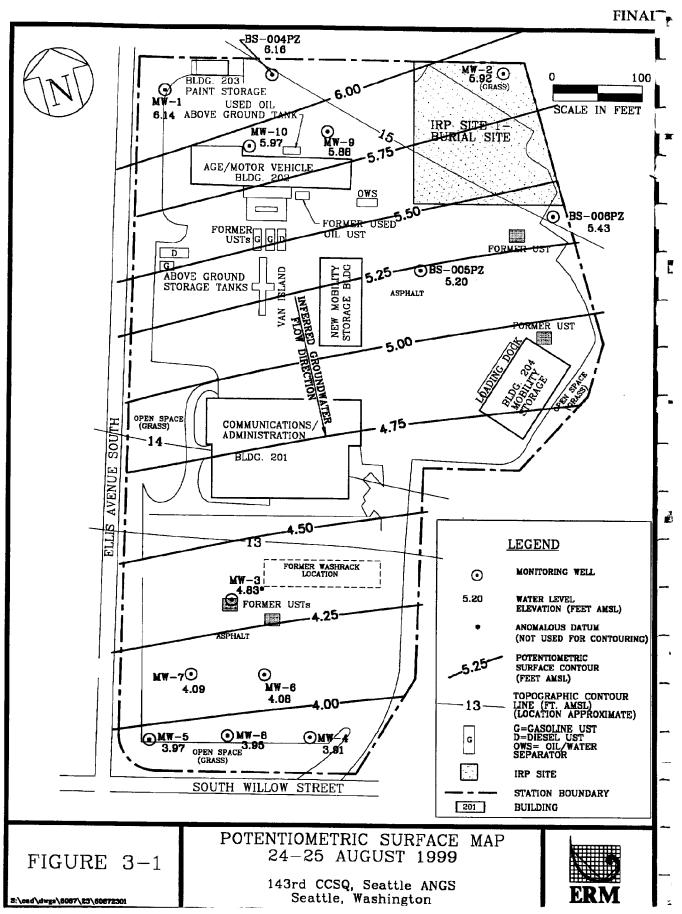
Static water levels measured in the monitoring wells on 24-25 August 1999 ranged from approximately 7 to 10 feet below ground surface. Depth measurements were converted to groundwater elevations by subtracting the measured depth to water in each well from the known elevation of the wellhead (top of well casing). A potentiometric surface map generated from the groundwater elevation data is presented on Figure 3-1. The inferred direction of groundwater flow was toward the south, generally consistent with previous measurements. Cumulative water level data for the Seattle ANGS monitoring wells are included in Appendix A.

3.2 Field Parameter Results

Before groundwater samples were collected at each monitoring well, the well was purged until field parameter measurements stabilized. The final field parameter measurements are summarized on Table 3-1.

3.3 Analytical Results

The analytical testing results for the August 1999 groundwater samples are summarized on Table 3-2. Table 3-2 also includes the results for groundwater samples collected during the RI and previous quarterly sampling events, and the PSGs derived from chemical-specific ARARs. The analytical testing results for the field QC blank samples are summarized on Table 3-3. Copies of the laboratory data summary



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TABLE 3-1
Final Field Parameter Measurements During Monitoring Well Purging, August 1999
143rd CCSQ, Seattle ANGS, Seattle, Washington

	***	a sa sa fil	elf Farameters		Alexander
Monitoribi W-II	Temperature (Degrees Celsius)	p) d	Specific Conductance (EX/cm)	Turbidity (NTU)	(Dissolved (Dxygen (mg/l)
BS-004PZ (Background Well)	15.8	5.59	285	19	6.18
BS-005PZ	16.0	5.52	288	15	0.27
BS-006PZ	15.9	5.54	340	10	0.34
MW-1 (Background Well)	16.2	5.64	322	19	0.89
MW-2	14.2	5.48	331	51	0.61
MW-3	18.9	5.60	338	10	0.43
MW-4	15.9	5.72	342	31	0.14
MW-5	16.4	5.61	288	15	0.20
MW-6	16.7	5.73	284	15	0.11
MW-7	16.4	5.69	281	32	0.18
MW-8	15.6	5.70	380	10	0.84
MW-9	15.6	5.31	290	53	0.28
MW-10	16.0	5.59	227	30	5.21

Notes:

pH = Acidity/alkalinity

μS/cm= Microsiemens per centimeter

NTU = Nephelometric turbidity units

mg/l = Milligrams per liter

TABLE 3-2

Organic Constituents Detected in Groundwater Monitoring Wells 143rd CCSQ, Seattle ANGS, Seattle, Washington

Location	Dalle j	Acetonie	e (A Chloroform	Toluëne	Bromodichloro methana	1,1- Clichlördethane	1,1,1 Trichloroethane	CB-1.2- Dichlorbethene	1,3,5 Trimethylberizene	Trichloroethene	Tetrachloroethene
	09/17/1996	ND	ND	(ND)	ND	0.3	3.7	ND	ND	ND	3.8
	9/17/96 (dup)	ND	ND	(ND)	ND	0.3	3.8	ND	ND	ND	3.8
	01/14/1997	ND	ND	ND	ND	ND	2.4	ND	ND	ND	5.1
	04/11/1997	ND	ND	ND	ND	ND	3.3	ND	ND	ND	17
BS-004PZ	07/10/1997	ND	ND	ND	ND	ND	1.8	ND	ND	ND	(ND)
(Background Well)	09/02/1998	ND	ND	ND	ND	ND	ND	ND	NA	ND	2.0
	11/25/1998	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	02/24/1999	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	05/19/1999	ND	ND	ND	ND	ND	ND	ND	NA	ND	6.8
	08/25/1999	ND	(ND)	ND	(ND)	ND	ND	ND	NA	ND	ND
	09/17/1996	ND	ND	ND	ND	ND	ND	ND	0.2	ND	ND
	01/15/1997	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	04/11/1997	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	07/11/1997	ND	ND	ND	ND	ND	ND	ND	ND	ND	4.7
BS-005PZ	09/01/1998	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	11/25/1998	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	02/24/1999	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	05/18/1999	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	08/24/1999	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	09/17/1996	ND	ND	ND	ND	ND	ND	ND	0.2	ND	ND
	01/14/1997	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	04/11/1997	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	07/11/1997	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
BS-006PZ	09/02/1998	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	11/24/1998	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	02/24/1999	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	05/18/1999	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	08/24/1999	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	10/18/1996	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	12/17/1996	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	01/14/1997	ND	ND	1.1	ND	ND	ND	ND	ND	ND	ND
	04/11/1997	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
MW-1	07/11/1997	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
(Background Well)	09/01/1998	ND	ND	6.0	ND	ND	ND	ND	NA	ND	ND
	11/25/1998	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	02/24/1999	ND	ND	ND	ND	ND	ND	ND	NA	ND	5.2
	05/19/1999	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	08/25/1999	ND	(ND)	ND	(ND)	ND	ND	ND	NA	ND	ND

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Organic Constituents Detected in Groundwater Monitoring Wells 143rd CCSQ, Seattle ANGS, Seattle, Washington

Location	Date	Acetype	Gilonoform	Toluene	Bromodichloro methane	Lil- Dichloroethane		Cis-L2- Dichigroethënë	1,3,5 Trimethylbenzene	Trichloroethene	Tetrachloroethene
	10/18/1 996	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	12/17/1996	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	01/15/1997	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	1/15/97 (dup)	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	04/10/1997	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
MW-2	07/11/1997	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	09/02/1998	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	11/25/1998	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	02/24/1999	Ð	ND	ND	ND	ND	ND	ND	NA	ND	ND
	05/18/1999	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	08/24/1999	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	10/18/1996	18	ND	ND	ND	ND	ND	ND	ND	ND	ND
	10/18/96 (dup)	20	ND	ND	ND	ND	ND	ND	ND	ND	ND
	12/17/1996	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	12/17/% (dup)	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	01/15/1997	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	04/11/1997	ND	ND	ND	ND	ND	ND	ND	ND	ŊD	ND
MW-3	07/11/1997	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	09/02/1998	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	9/2/98 (dup)	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	11/24/1998	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	11/24/98 (dup)	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	02/25/1999	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	05/18/1999	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	08/24/1999	ND	ND	ND	ND	ND	ND	ND	NA	ND	1.2
	10/18/1996	11	ND	ND	ND	ND	ND	ND	ND	3.9	ND
	12/17/1996	ND	ND	ND	ND	ND	ND	ND	ND	2.7	ND
	01/14/1997	ND	ND	ND	ND	ND	ND	ND	ND	3.4	ND
	04/11/1997	ND	ND	ND	ND	ND	ND	ND	ND	3.2	ND
	07/11/1997	ND	ND	ND	ND	ND	ND	ND	ND	2.8	ND
MW-4	09/02/1998	ND	ND	ND	ND	ND	ND	ND	NA	2.0	ND
	11/24/1998	ND	ND	ND	ND	ND	ND	ND	NA	3.4	ND
	02/24/1999	ND	ND	ND	ND	ND	ND	ND	NA	2.6	ND
	2/24/99 (dup)	ND	ND	ND	ND	ND	ND	ND	NA	2.6	ND
	05/18/1999	ND	ND	ND	ND	ND	ND	ND	NA	2.9	ND
	08/24/1999	ND	ND	ND	ND	ND	ND	ND	NA	3.3	ND

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TABLE 3-2

Organic Constituents Detected in Groundwater Monitoring Wells 143rd CCSQ, Seattle ANGS, Seattle, Washington

Location	Date	Acetone	Chloroform	Toluene	Bromoeichioro methane	Dichioroethane	Lile : Trichloroethane	Dichloroethene	1.15. Friedelingsbenzene	Trichloroethene	Tetrachloroethen
	10/18/1996	ND	ND	ND	ND	ND	ND	5.6	ND	ND	ND
	12/17/1996	ND	ND	ND	ND	ND	ND	4.9	ND	ND	ND
	01/14/1997	ND	ND	ND	ND	ND	ND	2.7	ND	ND	ND
	04/11/1997	ND	ND	ND	ND	ND	ND	1.4	ND	ND	ND
	4/11/97 (dup)	ND	ND	ND	ND	ND	ND	1.6	ND	ND	ND
MW-5	07/10/1997	ND	ND	ND	ND	ND	ND	3.5	ND	2.1	(ND)
	7/10/97 (dup)	ND	ND	ND	ND	ND	ND	2.8	ND	ND	ND
	09/02/1998	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	11/24/1998	ND	ND	ND	ND	ND	ND	3.2	NA	ND	ND
	02/25/1999	ND	ND	ND	ND	ND	ND	1.7	NA	ND	ND
	05/18/1999	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	08/24/1999	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	09/02/1998	ND	ND	ND	ND	ND	ND	ND	NA	3.0	ND
	11/24/1998	ND	ND	ND	ND	ND	ND	ND	NA	3.0	ND
MW-6	02/25/1999	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	05/18/1999	ND	ND	ND	ND	ND	ND	1.0	NA	5.7	ND
	08/24/1999	ND	ND	ND	ND	ND	ND	ND	NA	3.5 [ND
	09/02/1998	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	11/24/1998	ND	ND	ND	ND	ND	ND	ND	NA	3.8	ND
MW-7	02/25/1999	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	05/18/1999	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	08/24/1999	ND	ND	ND	ND	ND	ND	ND	NA NA	1.9]	1.7]
	09/02/1998	ND	ND	ND	ND	ND	ND	ND	NA	3.0	ND
	11/24/1998	ND	ND	ND	ND	ND	ND	ND	NA	3.3	ND
	02/24/1999	ND	ND	ND	ND	ND	ND	39	NA	83	ND
MW-8	2/24/99 (dup)	ND	ND	ND	ND	ND	ND	42	NA	**** 87	ND
	05/18/1999	ND	ND	ND	ND	ND	ND	4.5		19	ND
	5/18/99 (dup)	ND	ND	ND	ND	ND	ND	4.7	NA	B. 21 -	ND
	08/24/1999	ND	ND	ND	ND	ND	ND	ND	NA	7.2	ND
	8/24/99 (dup)	ND	ND	ND	ND	ND	ND	ND	NA	67 J	ND
	09/01/1998	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	11/24/1998	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
MW-9	02/24/1999	ND	ND	ND	ND	ND	ND	ND	NA NA	ND	ND
	05/18/1999	ND	ND	ND	ND	ND	1.4	ND	NA NA	ND	1.8
	08/24/1999	ND	ND	ND	ND	ND	2.0	ND	NA NA	ND	1.9

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Organic Constituents Detected in Groundwater Monitoring Wells 143rd CCSQ, Seattle ANGS, Seattle, Washington

Location :	Dale	Acadiste	Chlarotonn	Tomene	Bromodichlore metione	Gl-A. Dictionsettune	Teresismostisms		13.5 Trinsettylbenzene	Trichloroethene	Tetrachloroethene
	09/01/1998	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	11/25/1998	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
MW-10	02/24/1999	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	05/19/1999	ND	ND	ND	ND	ND	ND	ND	NA	ND	ND
	08/25/1999	ND	(ND)	ND	(ND)	ND	ND	ND	NA	ND	ND
RI Proje	ect Screening Goal	800	_	40	-	800	200	70	0.507	5.0	5.0

Notes:

All concentrations in micrograms per liter (µg/l).

ND = Not detected above laboratory method reporting limit.

(ND) = A positive detection was reported by the laboratory for this constituent in the sample indicated. The sample result was qualified as "not detected" based on a detection of the constituent in an associated quality control blank (United States Environmental Protection Agency Contract Laboratory Program "10x" and "5x" rules).

NA = Not analyzed

dup = Duplicate sample

] = Estimated concentration

RI = Remedial Investigation

Shaded cell/bold typeface indicates a value exceeding the associated RI project screening goal.

-- = RI project screening goal not established

3-/

			inic Compounds	
Secretary Type 2 1 2 per second	orn Date	Sample Number	Bromo- dichloromethane	Chloroform
Field Blank (Tap Water)	08/24/1999	MW-8-99/00-1FT	1.6	45
Field Blank (ASTM Type II Water)	08/24/1999	MW-8-99/00-1FA	ND	ND
Rinsate Blank	08/24/1999	MW-4-99/00-1R	ND	ND
Trip Blank	08/24/1999	TB082499-1	ND	ND

Notes:

All concentrations in micrograms per liter (µg/l).

ASTM = American Society for Testing and Materials

ND = Not detected above laboratory method reporting limit.

3-8

sheets are included in Appendix B. The QC Data Validation Report is included in Appendix C. Appendix D contains copies of the Chain-of-Custody Forms.

The only analyte detected above PSGs in August 1999 was TCE. The concentration of TCE in well MW-8 (7.2 μ g/l) exceeded the MTCA Method A Cleanup Level of 5.0 μ g/l for this compound. TCE also was detected below the PSG in the groundwater samples collected from wells MW-4, MW-6, and MW-7 in the southern portion of the Station. PCE was detected at concentrations below the PSG in the groundwater samples collected from wells MW-3, MW-7, and MW-9. 1,1,1-trichloroethane was detected below the PSG in the groundwater sample from well MW-9.

Field QC blank samples (Table 3-3) were analyzed for the same parameters as the associated project samples. Chloroform and bromodichloromethane were detected in the tap water field blank. These compounds are common residual byproducts of the drinking water chlorination process, and were likely present in the source water used to prepare the tap water field blank. Chloroform and bromodichloromethane were also detected in three groundwater samples. However, in accordance with the USEPA Contract Laboratory Program National Functional Guidelines for Organic Data Review (February 1994), these detections were qualified as "not detected" based on the presence of these compounds in the tap water field blank.

SECTION 4.0

CONCLUSIONS

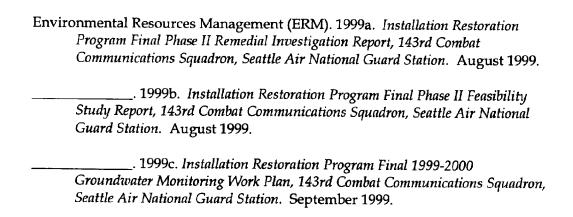
Groundwater samples collected at the Seattle ANGS in August 1999 were analyzed for VOCs. TCE was the only analyte detected above PSGs. TCE was detected above the Washington State MTCA Method A Cleanup Level of 5.0 μ g/l in one monitoring well (MW-8) near the southern Station boundary. The detected concentration (7.2 μ g/l) was less than the concentrations detected in this well in February 1999 (83 μ g/l) and May 1999 (19 μ g/l).

1

The source of the TCE in groundwater has not been identified. As discussed in the Phase II RI Report (ERM 1999a), the TCE may be related to the dissolved VOC plume in shallow groundwater beneath the Boeing facility immediately south of the Seattle ANGS.

SECTION 5.0

REFERENCES



APPENDIX A

WATER LEVEL DATA

TABLE A-1

Monitoring Well Water Level Summary

143rd CCSQ, Seattle ANGS, Seattle, Washington

Monitoring Well	Measuring Point Elevation (ft amsl)	Date	Depth to Water (ft bmp)	Water Level Elevation (ft amsl)
		9/17/96	8.88	5.78
*		10/22/96	8.93	5.73
		12/17/96	8.08	6.58
	14.66	1/14/97	6.98	7.68
BS-004PZ		4/11/97	7.23	7.43
(Background Well)		7/10/97	8.08	6.58
		9/1/98	9.79	4.87
		11/24/98	9.39	5.27
		2/24/99	6.84	7.82
	ļ	5/18/99	7.81	6.85
		8/25/99	8.50	6.16
		9/17/96	9.16	5.23
		10/22/96	9.42	4.97
		12/17/96	8.51	5.88
		1/15/97	7.48	6.91
	14.39	4/10/97	7.65	6.74
BS-005PZ		7/11/97	8.47	5.92
		9/1/98	10.12	4.27
		11/24/98	9.41	4.98
		2/24/99	7.32	7.07
		5/18/99	8.15	6.24
		8/24/99	9.19	5.20
		9/17/96	9.12	5.47
		10/22/96	9.47	5.12
		12/17/96	8.54	6.05
		1/14/97	7.62	6.97
		4/11/97	7.77	6.82
BS-006PZ	14.59	7/11/97	8.49	6.10
		9/1/98	10.29	4.30
		11/24/98	9.37	5.22
		2/24/99	7.42	7.17
		5/28/99	8.20	6.39
		8/24/99	9.16	5.43

FINAL

TABLE A-1

Monitoring Well Water Level Summary
143rd CCSQ, Seattle ANGS, Seattle, Washington

Monitoring Well	Measuring Point Elevation (ft amsl)	Date	Depth to Water (ft bmp)	Water Level Elevation (ft amsl)
MW-1 (Background Well)	14.92	10/22/96	9.18	5.74
		12/17/96	8.20	6.72
		1/14/97	7.11	7.81
		4/10/97	7.58	7.34
		7/11/97	8.51	6.41
		9/1/98	10.22	4.70
		11/24/98	9.45	5.47
		2/24/99	7.12	7.80
		5/18/99	8.25	6.67
		8/25/99	8.78	6.14
	14.60	10/22/96	8.89	5. <i>7</i> 1
		12/17/96	8.03	6.57
		1/15/97	7.13	7.47
		4/10/97	7.25	7.35
MW-2		7/11/97	7.98	6.62
		9/1/98	9.59	5.01
		11/24/98	9.75	4.85
		2/24/99	6.70	7.90
		5/18/99	7.71	6.89
		8/24/99	8.68	5.92
MW-3	11.88	10/22/96	7.77	4.11
		12/17/96	6.78	5.10
		1/15/97	7.80	4.08
		4/11/97	6.06	5.82
		7/11/97	6.94	4.94
		9/1/98	8.09	3.79
		11/24/98	7.20	4.68
		2/24/99	5.56	6.32
		5/18/99	6.65	5.23
		8/24/99	7.05	4.83
MW-4	12.05	10/22/96	8.20	3.85
		12/17/96	7.21	4.84
		1/14/97	6.31	5.74
		4/11/97	6.65	5.40
		7/11/97	7.43	4.62
		9/1/98	8.21	3.84
		11/24/98	8.14	3.91
		2/24/99	6.08	5.97
		5/18/99	7.16	4.89
		8/24/99	8.14	3.91

TABLE A-1

Monitoring Well Water Level Summary
143rd CCSQ, Seattle ANGS, Seattle, Washington

Monitoring Well	Measuring Point Elevation (ft amsl)	Date	Depth to Water (ft bmp)	Water Level Elevation (ft amsl)
MW-5	13.94	10/22/96	10.06	3.88
		12/17/96	9.06	4.88
		1/14/97	8.01	5.93
		4/11/97	8.36	5.58
		7/10/97	9.23	4.71
		9/1/98	10.15	3.79
		11/24/98	10.11	3.83
		2/24/99	7.84	6.10
		5/18/99	8.98	4.96
		8/24/99	9.97	3.97
	11.62	9/1/98	8.38	3.24
		11/24/98	7.64	3.98
MW-6		2/24/99	5.50	6.12
		5/18/99	6.55	5.07
		8/24/99	7.54	4.08
MW-7	12.17	9/1/98	6.75	5.42
		11/24/98	7.30	4.87
		2/24/99	5.94	6.23
		5/18/99	7.05	5.12
		8/24/99	8.08	4.09
MW-8	11.90	9/1/98	8.89	3.01
		11/24/98	8.02	3.88
		2/24/99	5.82	6.08
		5/18/99	6.95	4.95
		8/24/99	7.95	3.95
MW-9	14.30	9/1/98	9.78	4.52
		11/24/98	8.00	6.30
		2/24/99	6.76	7.54
		5/18/99	7.69	6.61
		8/24/99	8.42	5.88
MW-10	14.97	9/1/98	10.42	4.55
		11/24/98	9.69	5.28
		2/24/99	7.40	7.57
		5/18/99	8.43	6.54
		8/25/99	9.00	5.97

Notes:

ft amsl = Feet above mean sea level

ft bmp = Feet below measuring point

APPENDIX B

LABORATORY DATA SUMMARY SHEETS



September 13, 1999

ERM-West 915 118th Avenue S.E. Suite 130 Bellevue WA 98005

Attention: Rob Leet

6067.23 (AG Project Number : 6007.23

Project Name : Seattle ANGS

Dear Mr. Leet:

On August 25, 1999, MultiChem Analytical Services received 18 samples for analysis. The samples were analyzed with EPA methodology or equivalent methods as specified in the attached analytical schedule. The results, sample cross reference, and quality control data are enclosed.

Sincerely,

Tamara B. Jerome Senior Project Manager

TBJ/hal/trm

Enclosure

cc: Anita Quesada

ERM-West, Inc. - CA

Laurie Depies ERM-Sacramento, CA

MultiChem
ANALYTICAL SERVICES

SAMPLE CROSS REFERENCE SHEET

CLIENT : ERM-WEST PROJECT # : 6007.23

PROJECT NAME : SEATTLE ANGS

MAS #	CLIENT DESCRIPTION	DATE SAMPLED	MATRIX
908046-1	TB082499-1	00/04/00	
908046-2	MW-8-99/00-1FA	08/24/99	WATER
908046-3	MW-8-99/00-1FT	08/24/99	WATER
908046-4	BS-006PZ-99/00-1	08/24/99	WATER
908046-5	MW-2-99/00-1	08/24/99	WATER
908046-6		08/24/99	WATER
908046-7	BS-005PZ-99/00-1	08/24/99	WATER
908046-8	MW-9-99/00-1	08/24/99	WATER
908046-9	MW-3-99/00-1	08/24/99	WATER
	MW-4-99/00-1	08/24/99	WATER
908046-10	MW-4-99/00-1R	08/24/99	WATER
908046-11	MW-8-99/00-1	08/24/99	WATER
908046-12	MW-8-99/00-1D	08/24/99	WATER
908046-13	MW-6-99/00-1	08/24/99	WATER
008046-14	MW-7-99/00-1	08/24/99	WATER
08046-15	MW-5-99/00-1	08/24/99	WATER
08046-16	BS-004PZ-99/00-1	08/25/99	WATER
008046-17	MW-10-99/00-1	08/25/99	WATER
908046-18	MW-1-99/00-1	08/25/99	WATER

---- TOTALS ----

MATRIX # SAMPLES
----WATER 18

MAS STANDARD DISPOSAL PRACTICE

The samples from this project will be disposed of in thirty (30) days from the date of the report. If an extended storage period is required, please contact our sample control department before the scheduled disposal date.

KCSlip4 42539

ANALYTICAL SCHEDULE

CLIENT : ERM-WEST
PROJECT # : 6007.23
PROJECT NAME : SEATTLE ANGS

ANALYSIS	TECHNIQUE REFERENCE		LAB	
VOLATILE ORGANICS ANALYSIS	GCMS	EPA 8260B	R	

R = MAS - Renton
ANC = MAS - Anchorage
SUB = Subcontract

MultiChem
ANALYTICAL SERVICES

CASE NARRATIVE

CLIENT : ERM-WEST PROJECT # : 6007.23

PROJECT NAME : SEATTLE ANGS

CASE NARRATIVE: VOLATILE ORGANICS ANALYSIS

The following anomalies were associated with the preparation and/or analysis of the samples in this accession:

Per client request, initial calibration for all quantified compounds has been performed using average response factors. In all cases, the %RSD is <30%. In order to achieve this %RSD criterion, some of the high or low end calibration points may have been eliminated from the initial calibration. However, the initial calibration of all target compounds contains at least five consecutive points over the calibration range, with the low point at or below the reporting limit, and the high point defining the upper limit of the calibration range.

The percent recovery of the surrogate spiking compound 1,4-dichloroethane-d4 exceeded the current MultiChem recovery limits of 81-130% in the samples identified as 908046-10 (MW-4-99/00-1R), 908046-12 (MW-8-99/00-1D), 908046-13 (MW-6-99/00-1), 908046-14 (MW-7-99/00-1), 908046-15 (MW-5-99/00-1), and the matrix spike/matrix spike duplicate (MS/MSD) analyses performed on sample 908046-11 (MW-8-99/00-1). Since the anomalies indicated a potential high bias, and target compounds were not detected in samples 908046-10 (MW-4-99/00-1F) and 908046-15 (MW-5-99/00-1), they were reported without reanalysis. Since all MS/MSD recoveries and associated relative percent difference (RPD) criteria were within current MultiChem acceptance criteria, the MS and MSD analyses were not reanalyzed. Samples 908046-12 (MW-8-99/00-1D), 908046-13 (MW-6-99/00-1) and 908046-14 (MW-7-99/00-1), were reanalyzed one day outside of the method recommended 14 day holding time for preserved water samples. Since the recoveries of 1,4-dichloroethane-d4 were within control limits in these analyses, both were reported. All surrogate anomalies have been flagged with an "H" for reporting purposes on the appropriate Form 2. No further corrective action was performed.

All other associated quality assurance/quality control (QA/QC) parameters were within established MultiChem control limits.

KCSlip4 42541



VOLATILE ORGANICS ANALYSIS DATA SUMMARY

CLIENT : ERM-WEST DATE SAMPLED : N/A DATE RECEIVED : N/A PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS DATE EXTRACTED : N/A CLIENT I.D. : METHOD BLANK DATE ANALYZED : 09/02/99 SAMPLE MATRIX : WATER UNITS : ug/L EPA METHOD : 8260B DILUTION FACTOR: 1 COMPOUNDS RESULTS CHLOROMETHANE <5.0 VINYL CHLORIDE <1.0 BROMOMETHANE <1.0 CHLOROETHANE <1.0 TRICHLOROFLUOROMETHANE <1.0 ACETONE <10 1,1-DICHLOROETHENE <1.0 METHYLENE CHLORIDE <5.0 CARBON DISULFIDE <5.0 TRANS-1, 2-DICHLOROETHENE <1.0 1,1-DICHLOROETHANE <1.0 VINYL ACETATE <10 2-BUTANONE <10 CHLOROFORM CIS-1, 2-DICHLOROETHENE 1,1,1-TRICHLOROETHANE 1,2-DICHLOROETHANE CARBON TETRACHLORIDE BENZENE <1.0 1,2-DICHLOROPROPANE <1.0 TRICHLOROETHENE <1.0 BROMODICHLOROMETHANE <1.0 CIS-1, 3-DICHLOROPROPENE <3.0 4-METHYL-2-PENTANONE <10 TRANS-1, 3-DICHLOROPROPENE <3.0 1, 1, 2-TRICHLOROETHANE <1.0 TOLUENE <1.0 CHLORODIBROMOMETHANE <2.0 2-HEXANONE <10 TETRACHLOROETHENE <1.0 CHLOROBENZENE <1.0 ETHYLBENZENE <1.0 BROMOFORM <3.0

MultiChem ANALYTICAL SERVICES

CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : METHOD BLANK SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED : N/A DATE RECEIVED : N/A DATE EXTRACTED : N/A DATE ANALYZED : 09/02/99 UNITS : ug/L DILUTION FACTOR : 1	
COMPOUNDS	RESULTS	
(m+p)-XYLENE STYRENE O-XYLENE 1,1,2,2-TETRACHLOROETHANE 1,3-DICHLOROBENZENE 1,4-DICHLOROBENZENE P-ISOPROPYLTOLUENE 1,2-DICHLOROBENZENE	<1.0 <1.0	
SURROGATE PERCENT RECOVERY	LIMITS	_
1,2-DICHLOROETHANE-D4 TOLUENE-D8 BROMOFLUOROBENZENE	125 91 80 - 120 94 75 - 118	



_	CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : METHOD BLANK SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED : N/A DATE RECEIVED : N/A DATE EXTRACTED : N/A DATE ANALYZED : 09/02/99 UNITS : ug/L DILUTION FACTOR : 1
-	COMPOUNDS	RESULTS
	CHLOROMETHANE VINYL CHLORIDE BROMOMETHANE	<5.0 <1.0
	CHLOROETHANE TRICHLOROFLUOROMETHANE	<1.0
-	1,1-DICHLOROETHENE METHYLENE CHLORIDE CARBON DISULFIDE	<1.0 <5.0
_	CARBON DISULFIDE TRANS-1,2-DICHLOROETHENE 1,1-DICHLOROETHANE VINYL ACETATE 2-BUTANONE	<1.0 <1.0 <10 <10
	CHLOROFORM CIS-1,2-DICHLOROETHENE 1,1,1-TRICHLOROETHANE	· - ·
	1,2-DICHLOROETHANE CARBON TETRACHLORIDE BENZENE 1,2-DICHLOROPROPANE	
_	1,2-DICHLOROPROPANE TRICHLOROETHENE BROMODICHLOROMETHANE CIS-1,3-DICHLOROPROPENE 4-METHYL-2-PENTANONE	<1.0 <1.0 <1.0
_	4-METHYL-2-PENTANONE TRANS-1, 3-DICHLOROPROPENE 1,1,2-TRICHLOROETHANE	<3.0 <10 <3.0 <1.0
	TOLUENE CHLORODIBROMOMETHANE 2-HEXANONE	∠10
	TETRACHLOROETHENE CHLOROBENZENE ETHYLBENZENE	<1.0 <1.0 <1.0
	BROMOFORM	<3.0



CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : METHOD BLANK SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED DATE RECEIVED DATE EXTRACTE DATE ANALYZED UNITS DILUTION FACT	D : N/A ED : N/A D : 09/02/99 : ug/L
COMPOUNDS	RESULTS	
(m+p)-XYLENE STYRENE O-XYLENE 1,1,2,2-TETRACHLOROETHANE 1,3-DICHLOROBENZENE 1,4-DICHLOROBENZENE P-ISOPROPYLTOLUENE 1,2-DICHLOROBENZENE	<1.0 <1.0 <1.0	
SURROGATE PERCENT RECOVERY		LIMITS
1,2-DICHLOROETHANE-D4 TOLUENE-D8 BROMOFLUOROBENZENE	124 94 92	81 - 130 80 - 120 75 - 118



_	CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : METHOD BLANK SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED : N/A DATE RECEIVED : N/A DATE EXTRACTED : N/A DATE ANALYZED : 09/05/99 UNITS : ug/L DILUTION FACTOR : 1
_	COMPOUNDS	
_	CHLOROMETHANE VINYL CHLORIDE BROMOMETHANE	<5.0 <1.0
	CHLOROETHANE TRICHLOROFLUOROMETHANE ACETONE 1,1-DICHLOROETHENE	
_	1,1-DICHLOROETHENE METHYLENE CHLORIDE CARBON DISULFIDE TRANS-1,2-DICHLOROETHENE 1,1-DICHLOROETHANE VINYL ACETATE	<5.0 <5.0 <1.0 <1.0
~ "	Z-BUTANONE	<10
	CHLOROFORM CIS-1,2-DICHLOROETHENE 1,1,1-TRICHLOROETHANE 1,2-DICHLOROETHANE	<1.0 <1.0 <1.0
	1,2-DICHLOROETHANE CARBON TETRACHLORIDE BENZENE 1,2-DICHLOROPROPANE	<1.0 <1.0 <1.0
_	TRICHLOROETHENE BROMODICHLOROMETHANE CIS-1,3-DICHLOROPROPENE 4-METHYL-2-PENTANONE	
	4-METHYL-2-PENTANONE TRANS-1,3-DICHLOROPROPENE 1,1,2-TRICHLOROETHANE TOLUENE CHLORODIBROMOMETHANE	<10 <3.0 <1.0
-	Z-HEXANONE	<10
	TETRACHLOROETHENE CHLOROBENZENE ETHYLBENZENE	
—	BROMOFORM	<3.0



CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : METHOD BLANK SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED DATE RECEIVE DATE EXTRACT DATE ANALYZE UNITS DILUTION FAC	D : N/A ED : N/A D : 09/05/99 : ug/L
COMPOUNDS	RESULTS	
(m+p)-XYLENE STYRENE O-XYLENE 1,1,2,2-TETRACHLOROETHANE 1,3-DICHLOROBENZENE 1,4-DICHLOROBENZENE P-ISOPROPYLTOLUENE 1,2-DICHLOROBENZENE	<1.0 <1.0 <1.0 <1.0 <2.0 <2.0 <2.0 <2.0	
SURROGATE PERCENT RECOVERY		LIMITS
1,2-DICHLOROETHANE-D4 TOLUENE-D8 BROMOFLUOROBENZENE	124 92 89	81 - 130 80 - 120 75 - 118



_	CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : METHOD BLANK SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED : N/A DATE RECEIVED : N/A DATE EXTRACTED : N/A DATE ANALYZED : 09/08/99 UNITS : ug/L DILUTION FACTOR : 1
> {	COMPOUNDS	
1		
_	CHLOROMETHANE VINYL CHLORIDE BROMOMETHANE CHLOROETHANE	<5.0 <1.0 <1.0
	TRICULOROFI HODOMERITAND	<1.0
	TRICHLOROFLUOROMETHANE	<1.0
	ACETONE 1,1-DICHLOROETHENE METHYLENE CHLORIDE CARBON DISULFIDE TRANS-1,2-DICHLOROETHENE	<10
	METHYLENE CHLORIDE	<1.0 <5.0
<i>,</i>	CARBON DISULFIDE TRANS-1,2-DICHLOROETHENE 1,1-DICHLOROETHANE VINYL ACETATE 2 PHTANONE	<5.0
	TRANS-1, 2-DICHLOROETHENE	<1.0
	1,1-DICHLOROETHANE	<1.0
-	VINYL ACETATE	<10
	2-BUTANONE	<10
	CHLOROFORM	<1.0
_	CHLOROFORM CIS-1,2-DICHLOROETHENE 1,1,1-TRICHLOROETHANE	<1.0
	1,1,1-TRICHLOROETHANE	<1.0
1	1,2-DICHLOROETHANE	<1.0
~	CARBON TETRACHLORIDE	<1.0
•	1,2-DICHLOROETHANE CARBON TETRACHLORIDE BENZENE 1,2-DICHLOROPROPANE	<1.0
	TRICHLOROETHENE	<1.0
		<1.0
	CTS-1.3-DICHLOROPPOPME	<1.0
	BROMODICHLOROMETHANE CIS-1,3-DICHLOROPROPENE 4-METHYL-2-PENTANONE	<3.0
	TRANS-1, 3-DICHLOROPROPENE	<3.0
R	1.1.2-TRICHIOROETHANE	<1.0
1	TOLUENE	<1.0
	TRANS-1, 3-DICHLOROPROPENE 1, 1, 2-TRICHLOROETHANE TOLUENE CHLORODIBROMOMETHANE	<2.0
	7 — H H X A N// INIH	-10
	TETRACHLOROETHENE CHLOROBENZENE ETHYLBENZENE	<1.0
	CHLOROBENZENE	<1.0
	ETHYLBENZENE	<1.0
_	BROMOFORM	<3.0



CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : METHOD BLANK SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED DATE RECEIVED DATE EXTRACTE DATE ANALYZED UNITS DILUTION FACT	D : N/A DD : N/A D : 09/08/99 : ug/L
COMPOUNDS	RESULTS	
(m+p)-XYLENE STYRENE O-XYLENE 1,1,2,2-TETRACHLOROETHANE 1,3-DICHLOROBENZENE 1,4-DICHLOROBENZENE P-ISOPROPYLTOLUENE 1,2-DICHLOROBENZENE	<1.0 <1.0 <1.0	
SURROGATE PERCENT RECOVERY		LIMITS
1,2-DICHLOROETHANE-D4 TOLUENE-D8 BROMOFLUOROBENZENE	118 93 91	81 - 130 80 - 120 75 - 118



	CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : TB082499-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED : 08/24/99 DATE RECEIVED : 08/25/99 DATE EXTRACTED : N/A DATE ANALYZED : 09/02/99 UNITS : ug/L DILUTION FACTOR : 1
 I	COMPOUNDS	
ı		
	CHLOROMETHANE VINYL CHLORIDE BROMOMETHANE	<5.0 <1.0
	CHLOROETHANE	<1.0
_	TRICHLOROFLUOROMETHANE	<1.0
	ACETONE	<10
	ACETONE 1,1-DICHLOROETHENE METHYLENE CHLORIDE	<1.0
_	METHYLENE CHLORIDE	<5.0
•	CARBON DISULFIDE	<5.0
	1 1-DICHLODOFTUNKE	<1.0
	CARBON DISULFIDE TRANS-1,2-DICHLOROETHENE 1,1-DICHLOROETHANE VINYL ACETATE	<1.0 <10
	2-BUTANONE	<10
	CHLOROFORM	<1.0
٠.	CHLOROFORM CIS-1,2-DICHLOROETHENE 1,1,1-TRICHLOROETHANE	<1.0
	1,1,1-TRICHLOROETHANE	<1.0
i	1,2-DICHLOROETHANE	<1.0
	CARBON TETRACHLORIDE	<1.0
	1,2-DICHLOROETHANE CARBON TETRACHLORIDE BENZENE 1,2-DICHLOROPROPANE	<1.0
	1,2-DICHLOROPROPANE	<1.0
	TRICHLOROETHENE	/1 0
	BROMODICHLOROMETHANE CIS-1,3-DICHLOROPROPENE 4-METHYL-2-PENTANONE	<1.0
	4 MEMILY 2 PRIMITANCE	<3.0
	#PANS_1 2-DICH OPODDODDOD	<10
	1,1,2-TRICHLOROETHANE	<3.0
	TOLIENE	<1.0
1	TRANS-1,3-DICHLOROPROPENE 1,1,2-TRICHLOROETHANE TOLUENE CHLORODIBROMOMETHANE	<2.0
	2-HEXANONE	<10
_	TETRACHLOROETHENE	<1.0
	CHLOROBENZENE	<1.0
	TETRACHLOROETHENE CHLOROBENZENE ETHYLBENZENE	<1.0
	BROMOFORM	<3.0



CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : TB082499-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED : 08/24/99 DATE RECEIVED : 08/25/99 DATE EXTRACTED : N/A DATE ANALYZED : 09/02/99 UNITS : ug/L DILUTION FACTOR : 1	
COMPOUNDS	RESULTS	-
(m+p)-XYLENE STYRENE O-XYLENE 1,1,2,2-TETRACHLOROETHANE 1,3-DICHLOROBENZENE 1,4-DICHLOROBENZENE P-ISOPROPYLTOLUENE 1,2-DICHLOROBENZENE	<1.0 <1.0 <1.0 <1.0 <2.0 <2.0 <2.0 <2.0	-
SURROGATE PERCENT RECOVERY	LIMITS	
1,2-DICHLOROETHANE-D4 TOLUENE-D8 BROMOFLUOROBENZENE	128 81 - 130 94 80 - 120 98 75 - 118	_



erina mila	CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : MW-8-99/00-1FA SAMPLE MATRIX : WATER EPA METHOD : 8260B	
	COMPOUNDS	
	CHLOROMETHANE VINYL CHLORIDE BROMOMETHANE	<5.0 <1.0 <1.0
	CHLOROETHANE TRICHLOROFLUOROMETHANE ACETONE	<1.0 <1.0
	ACETONE 1,1-DICHLOROETHENE METHYLENE CHLORIDE CARBON DISULFIDE TRANS-1 2-DICHLOROETHENE	<1.0 <5.0 <5.0
_	CARBON DISULFIDE TRANS-1,2-DICHLOROETHENE 1,1-DICHLOROETHANE VINYL ACETATE 2-BUTANONE	<10
	CHLOROFORM CIS-1,2-DICHLOROETHENE 1,1,1-TRICHLOROETHANE	<1.0 <1.0 <1.0
	1,2-DICHLOROETHANE CARBON TETRACHLORIDE BENZENE 1,2-DICHLOROPROPANE	<1.0 <1.0 <1.0
-	TRICHLOROETHENE BROMODICHLOROMETHANE CIS-1,3-DICHLOROPROPENE 4-METHYL-2-PENTANONE	<1.0 <1.0 <1.0 <3.0
	TRANS-1, 3-DICHLOROPROPENE 1,1,2-TRICHLOROETHANE TOLUENE CHLORODIBROMOMETHANE	<3.0 <1.0 <1.0
	CHLORODIBROMOMETHANE 2-HEXANONE TETRACHLOROETHENE CHLOROBENZENE ETHYLBENZENE	
-2-	ETHYLBENZENE BROMOFORM	<1.0 <1.0 <3.0



CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : MW-8-99/00-1FA SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED : 08/24/99 DATE RECEIVED : 08/25/99 DATE EXTRACTED : N/A DATE ANALYZED : 09/02/99 UNITS : ug/L DILUTION FACTOR : 1
COMPOUNDS	RESULTS
(m+p)-XYLENE STYRENE O-XYLENE 1,1,2,2-TETRACHLOROETHANE 1,3-DICHLOROBENZENE 1,4-DICHLOROBENZENE P-ISOPROPYLTOLUENE 1,2-DICHLOROBENZENE	<1.0 <1.0
SURROGATE PERCENT RECOVERY	LIMITS
1,2-DICHLOROETHANE-D4 TOLUENE-D8 BROMOFLUOROBENZENE	127 81 - 130 94 80 - 120 100 75 - 118



_	CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : MW-8-99/00-1FT SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED DATE RECEIVED DATE EXTRACTE DATE ANALYZED UNITS DILUTION FACT	0 : 08/25/99 CD : N/A 0 : 09/02/99 : ug/L
	COMPOUNDS	RESULTS	
	(m+p)-XYLENE STYRENE O-XYLENE 1,1,2,2-TETRACHLOROETHANE 1,3-DICHLOROBENZENE 1,4-DICHLOROBENZENE P-ISOPROPYLTOLUENE 1,2-DICHLOROBENZENE	<1.0 <1.0 <1.0 <1.0 <2.0 <2.0 <2.0 <2.0	
-	SURROGATE PERCENT RECOVERY		LIMITS
مد	1,2-DICHLOROETHANE-D4 TOLUENE-D8BROMOFLUOROBENZENE	127 92 96	81 - 130 80 - 120 75 - 118

TRANS-1, 3-DICHLOROPROPENE

1,1,2-TRICHLOROETHANE

CHLORODIBROMOMETHANE

TETRACHLOROETHENE

TOLUENE

2-HEXANONE

CHLOROBENZENE

ETHYLBENZENE

BROMOFORM

CLIENT



VOLATILE ORGANICS ANALYSIS DATA SUMMARY

CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : BS-006PZ-99/00-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B	DILUTION FACTOR : 1
COMPOUNDS	RESULTS
CHLOROMETHANE VINYL CHLORIDE BROMOMETHANE CHLOROETHANE TRICHLOROFLUOROMETHANE ACETONE 1,1-DICHLOROETHENE METHYLENE CHLORIDE CARBON DISULFIDE TRANS-1,2-DICHLOROETHENE 1,1-DICHLOROETHANE VINYL ACETATE 2-BUTANONE CHLOROFORM CIS-1,2-DICHLOROETHENE 1,1,1-TRICHLOROETHANE 1,2-DICHLOROETHANE 1,2-DICHLOROETHANE 1,2-DICHLOROETHANE 1,2-DICHLOROFOPANE TRICHLOROETHENE BENZENE 1,2-DICHLOROPROPANE TRICHLOROETHENE BROMODICHLOROMETHANE CIS-1,3-DICHLOROMETHANE	<5.0 <1.0 <1.0 <1.0 <1.0 <1.0 <5.0 <5.0 <1.0 <1.0 <1.0 <1.0 <1.0 <1.0 <1.0 <1

<3.0

<1.0

<1.0

<2.0

<10

<1.0

<1.0

<1.0

<3.0

MultiChem
ANALYTICAL SERVICES

MAS I.D. # 908046-4

	CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : BS-006PZ-99/00-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED DATE RECEIVED DATE EXTRACTED DATE ANALYZED UNITS DILUTION FACTO	: 08/25/99 D : N/A : 09/02/99 : ug/L
-	COMPOUNDS	RESULTS	
	(m+p)-XYLENE STYRENE O-XYLENE 1,1,2,2-TETRACHLOROETHANE 1,3-DICHLOROBENZENE 1,4-DICHLOROBENZENE P-ISOPROPYLTOLUENE 1,2-DICHLOROBENZENE	<1.0 <1.0 <1.0	
	SURROGATE PERCENT RECOVERY		LIMITS
	1,2-DICHLOROETHANE-D4 TOLUENE-D8	127 94 95	81 - 130 80 - 120 75 - 118

2-BUTANONE

BROMOFORM



VOLATILE ORGANICS ANALYSIS DATA SUMMARY

CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : MW-2-99/00-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED : 08/24/99 DATE RECEIVED : 08/25/99 DATE EXTRACTED : N/A DATE ANALYZED : 09/02/99 UNITS : ug/L DILUTION FACTOR : 1
COMPOUNDS	RESULTS
CHLOROMETHANE VINYL CHLORIDE BROMOMETHANE CHLOROETHANE TRICHLOROFLUOROMETHANE ACETONE 1,1-DICHLOROETHENE METHYLENE CHLORIDE CARBON DISULFIDE TRANS-1,2-DICHLOROETHENE 1,1-DICHLOROETHANE VINYL ACETATE	<1.0 <10

<10

<10

<3.0

BROMODICHLOROMETHANE	VI.0
	<1.0
CIS-1, 3-DICHLOROPROPENE	
A MEMILIA O DESCRIPTION DIVI	<3.0
4-METHYL-2-PENTANONE	<10
TRANS-1, 3-DICHLOROPROPENE	
175 BICHBOROFROFENE	<3.0
1,1,2-TRICHLOROETHANE	<1.0
TOLUENE	<t.0< td=""></t.0<>
IOLOENE	<1.0
CHLORODIBROMOMETHANE	40.0

Z-HEXANONE	<10
TETRACHLOROETHENE	\10
	<1.0
CHLOROBENZENE	<1.0
ETHYLBENZENE	<1.U
	<1.0
RPOMOTORM	

KCSlip4 42557



	CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : MW-2-99/00-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED DATE RECEIVED DATE EXTRACTE DATE ANALYZED UNITS DILUTION FACT	0 : 08/25/99 CD : N/A 0 : 09/02/99 : ug/L
1	COMPOUNDS	RESULTS	
ł			
	O-XYLENE 1,1,2,2-TETRACHLOROETHANE 1,3-DICHLOROBENZENE 1,4-DICHLOROBENZENE P-ISOPROPYLTOLUENE 1,2-DICHLOROBENZENE	<1.0 <1.0	
_	SURROGATE PERCENT RECOVERY		LIMITS
	1,2-DICHLOROETHANE-D4	127	81 - 130
	TOLUENE-D8	90	80 - 120
_	BROMOFLUOROBENZENE	96	75 - 118

MultiChem ANALYTICAL SERVICES

Z-HEXANONE Z-HEXANONE TETRACHLOROETHENE CHLOROBENZENE ETHYLBENZENE STOROGORM C1.0 C1.0 C1.0 C1.0 C1.0	CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : BS-005PZ-99/00-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED : 08/24/99 DATE RECEIVED : 08/25/99 DATE EXTRACTED : N/A DATE ANALYZED : 09/02/99 UNITS : ug/L DILUTION FACTOR : 1	***
CHLOROMETHANE VINYL CHLORIDE SROMOMETHANE CHLOROETHANE CHLOROETHANE CHLOROFLUOROMETHANE ACETONE 1,1-DICHLOROETHENE 1,1-DICHLOROETHENE 1,1-DICHLOROETHENE CARBON DISULFIDE CARBON DISULFIDE VINYL ACETATE VINYL ACETATE VINYL ACETATE 1,0 CHLOROFTOM CIS-1,2-DICHLOROETHENE 1,1-DICHLOROETHANE CHLOROFORM CIS-1,2-DICHLOROETHANE CHLOROFORM CIS-1,2-DICHLOROETHANE CARBON TETRACHLORIDE CARBON TETRACHLOROPROPENE CIS-1,3-DICHLOROPROPENE CIS-1,0-DICHLOROPROPENE CIS-	COMPOUNDS	RESULTS	-
10.10	CHLOROMETHANE VINYL CHLORIDE BROMOMETHANE CHLOROETHANE TRICHLOROFLUOROMETHANE ACETONE 1,1-DICHLOROETHENE METHYLENE CHLORIDE CARBON DISULFIDE TRANS-1,2-DICHLOROETHENE 1,1-DICHLOROETHANE VINYL ACETATE 2-BUTANONE CHLOROFORM CIS-1,2-DICHLOROETHENE 1,1,1-TRICHLOROETHANE 1,2-DICHLOROETHANE CARBON TETRACHLORIDE BENZENE 1,2-DICHLOROPROPANE TRICHLOROETHENE BROMODICHLOROMETHANE CIS-1,3-DICHLOROPROPENE 4-METHYL-2-PENTANONE TRANS-1,3-DICHLOROPROPENE 1,1,2-TRICHLOROETHANE CHLOROBENE CHLORODIBROMOMETHANE CHLORODIBROMOMETHANE CHLORODIBROMOMETHANE CHLOROBENZENE ETHYLBENZENE ETHYLBENZENE ETHYLBENZENE	<5.0 <1.0 <1.0 <1.0 <1.0 <5.0 <5.0 <5.0 <5.0 <1.0 <1.0 <1.0 <1.0 <1.0 <1.0 <1.0 <1	-



	CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : BS-005PZ-99/00-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE EXTRACTE DATE ANALYZED UNITS DILUTION FACT	0 : 08/25/99 D : N/A 0 : 09/02/99 : ug/L
1	COMPOUNDS	RESULTS	
	(m+p)-XYLENE STYRENE O-XYLENE 1,1,2,2-TETRACHLOROETHANE 1,3-DICHLOROBENZENE 1,4-DICHLOROBENZENE P-ISOPROPYLTOLUENE 1,2-DICHLOROBENZENE	<1.0 <1.0 <1.0	
	SURROGATE PERCENT RECOVERY		LIMITS
- -	1,2-DICHLOROETHANE-D4 TOLUENE-D8	124 90 98	81 - 130 80 - 120 75 - 118

MultiChem ANALYTICAL SERVICES

PROJECT # PROJECT NAME CLIENT I.D. SAMPLE MATRIX EPA METHOD	: ERM-WEST : 6007.23 : SEATTLE ANGS : MW-9-99/00-1 : WATER : 8260B	DATE SAMPLED : 08/24/99 DATE RECEIVED : 08/25/99 DATE EXTRACTED : N/A DATE ANALYZED : 09/02/99 UNITS : ug/L DILUTION FACTOR : 1	•
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COMPOUNDS		_
CHLOROMETHANE	<5.0	
VINYL CHLORIDE BROMOMETHANE CHLOROETHANE TRICHLOROFILIOROMETHANE	<1.0	
BROMOMETHANE	<1.0	
CHLOROETHANE	<1.0	
TRICHLOROFLUOROMETHANE	<1.0	
ACETONE 1,1-DICHLOROETHENE METHYLENE CHLORIDE CARBON DISULFIDE TRANS-1,2-DICHLOROETHENE	<10	
1,1-DICHLOROETHENE	<1.0	
METHYLENE CHLORIDE	<5.0	•
CARBON DISULFIDE	<5.0	
TRANS-1, 2-DICHLOROETHENE	<1.0	
1,1-DICHLOROETHANE	<1.0	
CARBON DISULFIDE TRANS-1,2-DICHLOROETHENE 1,1-DICHLOROETHANE VINYL ACETATE 2-BUTANONE	<10	_
2 DOLLETOND	<10	4
CHLOROFORM	<1.0	
CIS-1,2-DICHLOROETHENE	<1.0 <1.0	
CHLOROFORM CIS-1,2-DICHLOROETHENE 1,1,1-TRICHLOROETHANE 1,2-DICHLOROETHANE	2.0	
1,2-DICHLOROETHANE CARBON TETRACHLORIDE BENZENE 1,2-DICHLOROPROPANE TRICHLOROETHENE	2.U <1.0	i
CARBON TETRACHLORIDE	<1.0 <1.0	
BENZENE	<1.0	-
1,2-DICHLOROPROPANE	<1.0	1
TRICHLOROETHENE	<1.0	
BROMODICHLOROMETHANE	<1.0	
TRICHLOROETHENE BROMODICHLOROMETHANE CIS-1, 3-DICHLOROPROPENE 4-METHYL-2-PENTANONE TRANS-1, 3-DICHLOROPROPENE	<1.0	_
4-METHYL-2-PENTANONE	<3.0	
TRANS-1, 3-DICHLOROPROPENE	<10	
1,1,2-TRICHLOROETHANE	<3.0	
TOLUENE	<1.0	-
TRANS-1, 3-DICHLOROPROPENE 1,1,2-TRICHLOROETHANE TOLUENE CHLORODIBROMOMETHANE 2-HEXANONE	<1.0	
2-HEXANONE	<2.0	
TETRACHLOROETHENE	<10	~
TETRACHLOROETHENE CHLOROBENZENE ETHYLBENZENE EROMOEODM	1.9	
ETHYLBENZENE	<1.0	
BROMOFORM		
	<3.0	



_	CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : MW-9-99/00-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED DATE RECEIVED DATE EXTRACTE DATE ANALYZED UNITS DILUTION FACT	: 08/25/99 D : N/A : 09/02/99 : ug/L
-	COMPOUNDS	RESULTS	
1			
	(m+p)-XYLENE STYRENE O-XYLENE 1,1,2,2-TETRACHLOROETHANE 1,3-DICHLOROBENZENE 1,4-DICHLOROBENZENE P-ISOPROPYLTOLUENE 1,2-DICHLOROBENZENE	<1.0 <1.0 <1.0 <1.0 <2.0 <2.0 <2.0 <2.0	
	SURROGATE PERCENT RECOVERY		LIMITS
_	1,2-DICHLOROETHANE-D4 TOLUENE-D8 BROMOFLUOROBENZENE	125 90 99	81 - 130 80 - 120 75 - 118



VOLATILE ORGANICS ANALYSIS DATA SUMMARY

EPA METHOD COMPOUNDS	: 8260B	UNITS : ug/L DILUTION FACTOR : 1
PROJECT # PROJECT NAME CLIENT I.D. SAMPLE MATRIX	: ERM-WEST : 6007.23 : SEATTLE ANGS : MW-3-99/00-1 : WATER	DATE SAMPLED : 08/24/99 DATE RECEIVED : 08/25/99 DATE EXTRACTED : N/A DATE ANALYZED : 09/02/99

COMPOUNDS	RESULTS
CHLOROMETHANE	
VINYI, CHIOPIDE	<5.0
BROMOMETHANE	<1.0
CHLOROFTHANE	<1.0
VINYL CHLORIDE BROMOMETHANE CHLOROETHANE TRICHLOROFLUOROMETHANE ACETONE	<1.0
ACETONE	<1.0
TRICHLOROFLUOROMETHANE ACETONE 1,1-DICHLOROETHENE METHYLENE CHLORIDE CARBON DISHLEIDE	<10
METHYLENE CHLORIDE	<1.0
CARBON DISHLEIDE	<5.0
TRANS-1 2-DICHI OPOETURATE	<5.0
1.1-DICHLOROFTUNKE	<1.0
TRANS-1, 2-DICHLOROETHENE 1, 1-DICHLOROETHANE VINYL ACETATE 2-BUTANONE	<1.0
2-BUTANONE	<10
CHIODOFORM	<10
CIS-1,2-DICHLOROETHENE 1,1,1-TRICHLOROETHANE 1,2-DICHLOROETHANE	<1.0
1.1.1-TRICHLOROFTUAME	<1.0
1, 2-DICHLOROETHANE	<1.0
CARBON TETRACHLORIDE	<1.0
1,2-DICHLOROETHANE CARBON TETRACHLORIDE BENZENE 1,2-DICHLOROPROPANE TRICHLOROETHENE	<1.0
1,2-DICHLOROPROPANE	<1.0
TRICHLOROETHENE	<1.0
BROMODICHLOROMETHANE CIS-1,3-DICHLOROPROPENE 4-METHYL-2-PENTANONE TRANS-1,3-DICHLOROPROPENE	<1.0
CIS-1.3-DICHLOPORPORENTE	<1.0
4-METHYL-2-DENTANONE	<3.0
TRANS-1.3-DICHLOPOPPOPPAR	<10
1.1.2-TRICHLOPOETUANTE	<3.0
TOLUENE	<1.0
TRANS-1,3-DICHLOROPROPENE 1,1,2-TRICHLOROETHANE TOLUENE CHLORODIBROMOMETHANE 2-HEXANONE	<1.0
2-HEXANONE	<2.0
TETRACHLOROETHENE	<10
CHLOROBENZENE	1.2
TETRACHLOROETHENE CHLOROBENZENE ETHYLBENZENE	<1.0
BROMOFORM	
DIVOLOT OLGI	<3.0

KCSlip4 42563



	CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : MW-3-99/00-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED DATE RECEIVED DATE EXTRACTE DATE ANALYZED UNITS DILUTION FACT	0 : 08/25/99 D : N/A 0 : 09/02/99 : ug/L
-	COMPOUNDS	RESULTS	
_ _	(m+p)-XYLENE STYRENE O-XYLENE 1,1,2,2-TETRACHLOROETHANE 1,3-DICHLOROBENZENE 1,4-DICHLOROBENZENE P-ISOPROPYLTOLUENE 1,2-DICHLOROBENZENE	<1.0 <1.0 <1.0 <1.0 <2.0 <2.0 <2.0 <2.0	
	SURROGATE PERCENT RECOVERY		LIMITS
	1,2-DICHLOROETHANE-D4 TOLUENE-D8BROMOFLUOROBENZENE	126 93 97	81 - 130 80 - 120 75 - 118

CIS-1,3-DICHLOROPROPENE

1,1,2-TRICHLOROETHANE

CHLORODIBROMOMETHANE

TETRACHLOROETHENE

CHLOROBENZENE

ETHYLBENZENE

TOLUENE

2-HEXANONE

BROMOFORM

TRANS-1, 3-DICHLOROPROPENE

4-METHYL-2-PENTANONE

: ERM-WEST

CLIENT



: 08/24/99

DATE SAMPLED

VOLATILE ORGANICS ANALYSIS DATA SUMMARY

PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : MW-4-99/00-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED : 08/24/99 DATE RECEIVED : 08/25/99 DATE EXTRACTED : N/A DATE ANALYZED : 09/05/99 UNITS : ug/L DILUTION FACTOR : 1	
COMPOUNDS	RESULTS	
CHLOROMETHANE VINYL CHLORIDE BROMOMETHANE CHLOROETHANE TRICHLOROFLUOROMETHANE ACETONE 1,1-DICHLOROETHENE METHYLENE CHLORIDE CARBON DISULFIDE TRANS-1,2-DICHLOROETHENE 1,1-DICHLOROETHANE VINYL ACETATE 2-BUTANONE CHLOROFORM CIS-1,2-DICHLOROETHENE 1,1,1-TRICHLOROETHANE 1,2-DICHLOROETHANE 1,2-DICHLOROETHANE CARBON TETRACHLORIDE BENZENE 1,2-DICHLOROPROPANE TRICHLOROETHENE	<5.0 <1.0 <1.0 <1.0 <1.0 <10 <1.0 <5.0 <5.0 <1.0 <1.0 <1.0 <1.0 <1.0 <1.0	
BROMODICHLOROMETHANE	<1.0	

<3.0

<10

<3.0

<1.0

<1.0

<2.0

<10

<1.0

<1.0

<1.0

<3.0



_	CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : MW-4-99/00-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED DATE RECEIVED DATE EXTRACTE DATE ANALYZED UNITS DILUTION FACT	0 : 08/25/99 D : N/A 0 : 09/05/99 : ug/L
	COMPOUNDS	RESULTS	
!			
	(m+p)-XYLENE STYRENE O-XYLENE 1,1,2,2-TETRACHLOROETHANE 1,3-DICHLOROBENZENE 1,4-DICHLOROBENZENE P-ISOPROPYLTOLUENE 1,2-DICHLOROBENZENE	<1.0 <1.0 <1.0 <1.0 <2.0 <2.0 <2.0 <2.0	
	SURROGATE PERCENT RECOVERY		LIMITS
^	1,2-DICHLOROETHANE-D4 TOLUENE-D8	129 93 92	81 - 130 80 - 120 75 - 118



VOLATILE ORGANICS ANALYSIS DATA SUMMARY

CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : MW-4-99/00-1R SAMPLE MATRIX : WATER EPA METHOD : 8260B COMPOUNDS	DATE SAMPLED : 08/24/99 DATE RECEIVED : 08/25/99 DATE EXTRACTED : N/A DATE ANALYZED : 09/05/99 UNITS : ug/L DILUTION FACTOR : 1	-
COMPOUNDS	RESULTS	_
BROMOMETHANE CHLOROETHANE	<5.0 <1.0 <1.0 <1.0	
TRICHLOROFLUOROMETHANE ACETONE	<1.0	
ACETONE 1,1-DICHLOROETHENE METHYLENE CHLORIDE CARBON DISULFIDE TRANS-1,2-DICHLOROETHENE 1,1-DICHLOROETHANE VINYL ACETATE 2-BUTANONE	<1.0 <1.0	
CARBON DISHLETDE	<5.0	_
TRANS-1,2-DICHLOROETHENE	<5.0 <1.0	
1,1-DICHLOROETHANE	<1.0	_
2-BUTANONE	<10	
CHLOROFORM CIS-1,2-DICHLOROETHENE 1,1,1-TRICHLOROETHANE 1,2-DICHLOROETHANE	<1.0	^
1,2-DICHLOROETHANE 1,2-DICHLOROETHANE CARBON TETRACHLORIDE BENZENE 1,2-DICHLOROPROPANE TRICHLOROETHENE BROMODICHLOROMETHANE	<1.0	
CARBON TETRACHLORIDE	<1.0	_
1 2-DICHI OPODRODANE	<1.0	
TRICHLOROETHENE	<1.0	
BROMODICHLOROMETHANE CIS-1,3-DICHLOROPROPENE 4-METHYL-2-PENTANONE TRANS-1,3-DICHLOROPROPENE	<1.0	_
CIS-1, 3-DICHLOROPROPENE	<3.0	
TRANS-1, 3-DICHLOROPROPENE	<10	
1,1,2-TRICHLOROETHANE	<3.0 <1.0	-
TRANS-1, 3-DICHLOROPROPENE 1,1,2-TRICHLOROETHANE TOLUENE CHLORODIBROMOMETHANE 2-HEXANONE	<1.0	
2-HEXANONE	<2.0	
TETRACHLOROETHENE	<10	
Z-HEXANONE TETRACHLOROETHENE CHLOROBENZENE ETHYLBENZENE BROMOFORM	<1.0 <1.0 <1.0	_
BROMOFORM	<3.0	

KCSlip4 42567



VOLATILE ORGANICS ANALYSIS DATA SUMMARY

CLIENT PROJECT # PROJECT NAME CLIENT I.D. SAMPLE MATRIX EPA METHOD	: ERM-WEST : 6007.23 : SEATTLE ANGS : MW-4-99/00-1R : WATER : 8260B	DATE EXTRACTED	: 08/24/99 : 08/25/99 : N/A : 09/05/99 : ug/L : 1
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-	COMPOUNDS	RESULTS	
	(m+p)-XYLENE STYRENE O-XYLENE 1,1,2,2-TETRACHLOROETHANE 1,3-DICHLOROBENZENE 1,4-DICHLOROBENZENE P-ISOPROPYLTOLUENE 1,2-DICHLOROBENZENE	<1.0 <1.0 <1.0 <1.0 <2.0 <2.0 <2.0 <2.0	
	SURROGATE PERCENT RECOVERY		LIMITS
_	1,2-DICHLOROETHANE-D4 TOLUENE-D8 BROMOFLUOROBENZENE	133 H 93 90	81 - 130 80 - 120 75 - 118

H = Out of limits.

TETRACHLOROETHENE

CHLOROBENZENE

ETHYLBENZENE

BROMOFORM



VOLATILE ORGANICS ANALYSIS DATA SUMMARY

		-
COMPOUNDS	5 T	
	RESULTS	
CHLOROMETHANE	<5.0	_
VINIL CHLORIDE	41 0	
BROMOMETHANE	<1.0	
BROMOMETHANE CHLOROETHANE TRICHLOROFLUOROMETHANE ACETONE	<1.0	
TRICHLOROFLUOROMETHANE	<1.0	_
ACETONE	<10	!
ACETONE 1,1-DICHLOROETHENE METHYLENE CHLORIDE CARBON DISH FIDE	<1.0	
METHILENE CHLORIDE	<5.0	
CARBON DISULFIDE	<5.0	
1 1 DICHLOROETHENE	<1.0	*
VINVI ACETATE	<1.0	
CARBON DISULFIDE TRANS-1,2-DICHLOROETHENE 1,1-DICHLOROETHANE VINYL ACETATE 2-BUTANONE	<10	
CHLOROFORM CIS-1,2-DICHLOROETHENE 1,1,1-TRICHLOROETHANE 1,2-DICHLOROETHANE	<1.0	
1.1.1-TRICHLOPOFTURNE	<1.0	~
1.2-DICHLOROETHANE	<1.0	
CARBON TETRACHLORIDE	<1.0	=
BENZENE	<1.0	
1,2-DICHLOROETHANE CARBON TETRACHLORIDE BENZENE 1,2-DICHLOROPROPANE TRICHLOROFTHENE	<1.0	
TRICHLOROETHENE	<1.0	
	7.2	
BROMODICHLOROMETHANE CIS-1,3-DICHLOROPROPENE 4-METHYL-2-PENTANONE TRANS-1 3-DICHLOROPROPENE	<1.0 <3.0	~
4-METHYL-2-PENTANONE	<10	1
TRANS-1, 3-DICHLOROPROPENE	<3.0	
1,1,2-TRICHLOROETHANE	<1.0	_
TOLUENE	<1.0	
TRANS-1, 3-DICHLOROPROPENE 1, 1, 2-TRICHLOROETHANE TOLUENE CHLORODIBROMOMETHANE 2-HEYANONE	<2.0	,
2 IIBIUUONE	<10	-
TETRACHLOROETHENE	11.0	

<1.0

<1.0

<1.0

<3.0



_	CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : MW-8-99/00-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED DATE RECEIVED DATE EXTRACTE DATE ANALYZED UNITS DILUTION FACT	: 08/25/99 D : N/A : 09/05/99 : ug/L
-	COMPOUNDS	RESULTS	
	(m+p)-XYLENE STYRENE O-XYLENE 1,1,2,2-TETRACHLOROETHANE 1,3-DICHLOROBENZENE 1,4-DICHLOROBENZENE P-ISOPROPYLTOLUENE 1,2-DICHLOROBENZENE	<1.0 <1.0 <1.0 <1.0 <2.0 <2.0 <2.0 <2.0	
_	SURROGATE PERCENT RECOVERY		LIMITS
بمر	1,2-DICHLOROETHANE-D4 TOLUENE-D8 BROMOFLUOROBENZENE	130 91 91	81 - 130 80 - 120 75 - 118



CLIENT PROJECT # PROJECT NAME CLIENT I.D. SAMPLE MATRIX EPA METHOD	: MW-8-99/00-1D	DATE SAMPLED : 08/24/99 DATE RECEIVED : 08/25/99 DATE EXTRACTED : N/A DATE ANALYZED : 09/06/99 UNITS : ug/L DILUTION FACTOR : 1
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COMPOUNDS	RESULTS
CHLOROMETHANE	
VINYL CHLORIDE	<5.0
BROMOME THANE	<1.0
CHLOROFTHAND	<1.0
TRICHLOROFLUOROMETHANE	
ACETONE	<1.0
ACETONE 1,1-DICHLOROETHENE METHYLENE CHLORIDE CARBON DISHLEDE	<10
METHYLENE CHLORIDE	<1.0
CARBON DISULFIDE	<5.0
TRANS-1.2-DICHLOROFTHENE	<5.0
CARBON DISULFIDE TRANS-1,2-DICHLOROETHENE 1,1-DICHLOROETHANE VINYL ACETATE 2-BUTANONE	<1.0
VINYL ACETATE	<1.0
2-BUTANONE	<10
CHI OBOEODM	<10
CILOROFORM CIS-1,2-DICHLOROETHENE 1,1,1-TRICHLOROETHANE	<1.0
1,1,1-TRICHLOROETHANE	<1.0
1,2-DICHLOROETHANE	<1.0
CARBON TETRACHLORIDE	<1.0
1,2-DICHLOROETHANE CARBON TETRACHLORIDE BENZENE 1,2-DICHLOROPROPANE TRICHLOROETHENE	<1.0
1,2-DICHLOROPROPANE	<1.0
TRICHLOROETHENE	<1.0
BROMODICHLOROMETHANE	6.7
CTS-1.3-DICHLOROPPOPENE	<1.0
BROMODICHLOROMETHANE CIS-1,3-DICHLOROPROPENE 4-METHYL-2-PENTANONE	<3.0
PRANS-1. 3-DICHLOPOPPOPENE	<10
1.1.2-TRICHLOPOETURNE	<3.0
OLUENE	<1.0
TRANS-1,3-DICHLOROPROPENE L,1,2-TRICHLOROETHANE TOLUENE CHLORODIBROMOMETHANE P-HEYANOME	<1.0
C-HEXANONE	<2.0
110121 0110	/10
ETRACHLOROETHENE CHLOROBENZENE CTHYLBENZENE	<1.0
THYLBENZENE	<1.0
ROMOFORM	<1.0
DIVITOR ONT	<3.0



VOLATILE ORGANICS ANALYSIS DATA SUMMARY

-	CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : MW-8-99/00-1D SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED DATE RECEIVED DATE EXTRACTE DATE ANALYZED UNITS DILUTION FACT	: 08/25/99 D : N/A : 09/06/99 : ug/L
)	COMPOUNDS	RESULTS	
_	(m+p)-XYLENE STYRENE O-XYLENE 1,1,2,2-TETRACHLOROETHANE 1,3-DICHLOROBENZENE 1,4-DICHLOROBENZENE P-ISOPROPYLTOLUENE 1,2-DICHLOROBENZENE	<1.0 <1.0	
	SURROGATE PERCENT RECOVERY		LIMITS
_	1,2-DICHLOROETHANE-D4 TOLUENE-D8 BROMOFLUOROBENZENE	136 H 90 92	81 - 130 80 - 120 75 - 118

H = Out of limits.

MAS I.D. # 908046-12

MAS I.D. # 908046-12RE

MultiChem ANALYTICAL SERVICES

VOLATILE ORGANICS ANALYSIS DATA SUMMARY

CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : MW-8-99/00-1D SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED : 08/24/99 DATE RECEIVED : 08/25/99 DATE EXTRACTED : N/A DATE ANALYZED : 09/08/99* UNITS : ug/L DILUTION FACTOR : 1	4
COMPOUNDS	RESULTS	•
CHLOROMETHANE	<5.0	_
VINYL CHLORIDE	<1.0	
BROMOMETHANE	<1.0	
CHLOROETHANE TRICHLOROETHANE	<1.0	
INICILLOROF LOOKOME THANE	<1.0	•
ACETONE		
1,1-DICHLOROETHENE	<1.0	
1,1-DICHLOROETHENE METHYLENE CHLORIDE CARBON DISH FIDE	<5.0	
CARBON DISULFIDE	<5.0	
CARBON DISULFIDE TRANS-1,2-DICHLOROETHENE 1,1-DICHLOROETHANE VINYL ACETATE 2-BUTANONE	<1.0	
1,1-DICHLOROETHANE	<1.0	
VINYL ACETATE	<10	•
	<10	
CHLOROFORM	· - ·	
CIS-1, 2-DICHLOROETHENE 1,1,1-TRICHLOROETHANE 1,2-DICHLOROETHANE	<1.0	_
1,1,1-TRICHLOROETHANE	<1.0	
1) 2 DICHLOROEIRANE	<1.0	
CARBON TETRACHLORIDE	<1.0	
BENZENE		•
1, 2-DICHLOROPROPANE	<1.0	
TRICHLOROETHENE	6.4	
BROMODICHLOROMETHANE		-
CIS-1, 3-DICHLOROPROPENE	<3.0	
CIS-1,3-DICHLOROPROPENE 4-METHYL-2-PENTANONE	<10	
TVANS-T, S-DICHHOROPROPENE	<3.0	
1,1,2 TRICHLOROETHANE	<1.0	-
I CALICARIA.	ITI	
CHLORODIBROMOMETHANE	<2.0	
Z-REARIONE	<10	_
TETRACHLOROETHENE		-
CHLOROBENZENE ETHYLBENZENE PROMOFORM	<1.0	
ETHYLBENZENE	<1.0	
BROMOFORM	<3.0	-
RE = Reanalyzed.		_

RE = Reanalyzed.
* = Reanalyzed outside of the recommended holding time.

MAS I.D. # 908046-12RE



VOLATILE ORGANICS ANALYSIS DATA SUMMARY

-	CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : MW-8-99/00-1D SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED : 08/24/99 DATE RECEIVED : 08/25/99 DATE EXTRACTED : N/A DATE ANALYZED : 09/08/99* UNITS : ug/L DILUTION FACTOR : 1
-	COMPOUNDS	RESULTS
_	(m+p)-XYLENE STYRENE O-XYLENE 1,1,2,2-TETRACHLOROETHANE 1,3-DICHLOROBENZENE 1,4-DICHLOROBENZENE P-ISOPROPYLTOLUENE 1,2-DICHLOROBENZENE	<1.0 <1.0 <1.0 <1.0 <2.0 <2.0 <2.0 <2.0

SURROGATE PERCENT RECOVERY LIMITS 1,2-DICHLOROETHANE-D4 81 - 130 80 - 120 75 - 118 119 TOLUENE-D8 92 BROMOFLUOROBENZENE 95

RE = Reanalyzed.

^{* =} Reanalyzed outside of the recommended holding time.



VOLATILE ORGANICS ANALYSIS DATA SUMMARY

CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : MW-6-99/00-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED : 08/24/99 DATE RECEIVED : 08/25/99 DATE EXTRACTED : N/A DATE ANALYZED : 09/06/99 UNITS : ug/L DILUTION FACTOR : 1	
COMPOUNDS	RESULTS	
CHLOROMETHANE VINYL CHLORIDE BROMOMETHANE CHLOROETHANE TRICHLOROFLUOROMETHANE ACETONE 1,1-DICHLOROETHENE METHYLENE CHLORIDE CARBON DISULFIDE	<5.0 <1.0 <1.0 <1.0 <1.0 <1.0 <1.0 <5.0 <5.0 <5.0 <1.0 <1.0 <1.0 <1.0 <1.0 <1.0 <1.0 <1	

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81 - 130 80 - 120 75 - 118

137 н

87

89

VOLATILE ORGANICS ANALYSIS DATA SUMMARY

_	CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : MW-6-99/00-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED : 08/24/99 DATE RECEIVED : 08/25/99 DATE EXTRACTED : N/A DATE ANALYZED : 09/06/99 UNITS : ug/L DILUTION FACTOR : 1
	COMPOUNDS	RESULTS
_	(m+p)-XYLENESTYRENE	<1.0 <1.0
	O-XYLENE 1,1,2,2-TETRACHLOROETHANE	<1.0 <1.0
	1,3-DICHLOROBENZENE	<2.0
	1,4-DICHLOROBENZENE P-ISOPROPYLTOLUENE	<2.0 <2.0
	1,2-DICHLOROBENZENE	<2.0
	SURROGATE PERCENT RECOVERY	LIMITS

H = Out of limits.

BROMOFLUOROBENZENE

TOLUENE-D8

1,2-DICHLOROETHANE-D4

MAS I.D. # 908046-13RE

CLIENT



CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : MW-6-99/00-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B COMPOUNDS	DATE SAMPLED : 08/24/99 DATE RECEIVED : 08/25/99 DATE EXTRACTED : N/A DATE ANALYZED : 09/08/99* UNITS : ug/L DILUTION FACTOR : 1
COMPOUNDS	RESULTS
CILODOMBETANT	
VINVI. CHI OD I DE	<5.0
BEOMOMETRANE	<1.0
VINYL CHLORIDE BROMOMETHANE CHLOROETHANE TRICHLOROETHOROMETHAND	<1.0
TRICHLOROFLUOROMETHANE	<1.0
ACETONE ACETONE	<1.0
1.1-DICHLOROFTHENE	<10
METHYLENE CHLORIDE	<1.0 <5.0
ACETONE 1,1-DICHLOROETHENE METHYLENE CHLORIDE CARBON DISULFIDE TRANS-1 2-DICHLOROFTHENE	<5.0
TRANS-1, 2-DICHLOROETHENE	<5.0
1,1-DICHLOROETHANE	<1.0
CARBON DISULFIDE TRANS-1,2-DICHLOROETHENE 1,1-DICHLOROETHANE VINYL ACETATE 2-BUTANONE	<1.0
2-BUTANONE	<10
CHLOROFORM	<1.0
CHLOROFORM CIS-1,2-DICHLOROETHENE 1,1,1-TRICHLOROETHANE 1,2-DICHLOROETHANE	<1.0
1,1,1-TRICHLOROETHANE	<1.0
1,2-DICHLOROETHANE	<1.0
CARBON TETRACHLORIDE	<1.0
BENZENE	<1.0
1,2-DICHLOROETHANE CARBON TETRACHLORIDE BENZENE 1,2-DICHLOROPROPANE TRICHLOROETHENE	<1.0
THE CONTRACT OF THE CONTRACT O	3.5
BROMODICHLOROMETHANE	<1.0
BROMODICHLOROMETHANE CIS-1,3-DICHLOROPROPENE 4-METHYL-2-PENTANONE	<3.0
4-METHIL-Z-PENTANONE	<10
1 1 2-TPICHLOROPROPENE	<3.0
TOLIENE	<1.0
CHLORODIRROMOMETURNE	<1.0
TRANS-1, 3-DICHLOROPROPENE 1,1,2-TRICHLOROETHANE TOLUENE CHLORODIBROMOMETHANE 2-HEXANONE	<2.0
TETRACHLOROETHENE CHLOROBENZENE ETHYLBENZENE BROMOFORM	<10
CHLOROBENZENE	<1.0
ETHYLBENZENE	<1.0
BROMOFORM	<1.0
	<3.0

RE = Reanalyzed.
* = Reanalyzed outside of the recommended holding time.



MAS I.D. # 908046-13RE

 CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : MW-6-99/00-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED : 08/24/99 DATE RECEIVED : 08/25/99 DATE EXTRACTED : N/A DATE ANALYZED : 09/08/99* UNITS : ug/L DILUTION FACTOR : 1
 COMPOUNDS	RESULTS
 (m+p)-XYLENE STYRENE O-XYLENE 1,1,2,2-TETRACHLOROETHANE 1,3-DICHLOROBENZENE 1,4-DICHLOROBENZENE P-ISOPROPYLTOLUENE 1,2-DICHLOROBENZENE	<1.0 <1.0 <1.0 <1.0 <2.0 <2.0 <2.0 <2.0

SURROGATE PERCENT RECOVERY		LIMITS
1,2-DICHLOROETHANE-D4 TOLUENE-D8 BROMOFLUOROBENZENE	123 90 93	81 - 130 80 - 120 75 - 118

RE = Reanalyzed.
* = Reanalyzed outside of the recommended holding time.



CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : MW-7-99/00-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED : 08/24/99 DATE RECEIVED : 08/25/99 DATE EXTRACTED : N/A DATE ANALYZED : 09/06/99 UNITS : ug/L DILUTION FACTOR : 1	_
COMPOUNDS	RESULTS	
CHLOROMETHANE	<5.0	
VINYL CHLORIDE	<1.0	
BROMOMETHANE		
CHLOROETHANE TRICHLOROFLUOROMETHANE	<1.0	
TRICHLOROFLUOROMETHANE	<1.0	_
ACETONE 1,1-DICHLOROETHENE METHYLENE CHLORIDE CAPRON DISH FIDE	<10	
1,1-DICHLOROETHENE	<1.0	
METHYLENE CHLORIDE	<5.0	-
CARBON DISULFIDE TRANS-1,2-DICHLOROETHENE 1,1-DICHLOROETHANE VINYL ACETATE 2-BUTTONE	<5.0	
1 AND CHI OPPERATE	<1.0	
T, T-DICHLOROETHANE	<1.0	
2-BUTANONE	<10	
5 DO ITMONE	/10	
CIS-1, 2-DICHLOPORTURNE	<1.0	
CHLOROFORM CIS-1,2-DICHLOROETHENE 1,1,1-TRICHLOROETHANE 1,2-DICHLOROETHANE	<1.0	-
CARBON TETRACHLORIDE	<1.0	
CARBON TETRACHLORIDE BENZENE 1,2-DICHLOROPROPANE TRICHLOROFTHENE	<1.0	_
1,2-DICHLOROPROPANE	<1.0	
INICALORUE I FINNE.	1 0	
BROMODICHLOROMETHANE	<1.0	_
BROMODICHLOROMETHANE CIS-1,3-DICHLOROPROPENE 4-METHYL-2-PENTANONE TRANS-1,3-DICHLOROPROPENE	<3.0	
4-METHYL-2-PENTANONE	<10	
TRANS-1, 3-DICHLOROPROPENE	<3 0	
1,1,2-TRICHLOROETHANE	<1 0	_
TOLUENE	<1.0	
TRANS-1, 3-DICHLOROPROPENE 1,1,2-TRICHLOROETHANE TOLUENE CHLORODIBROMOMETHANE 2-HEXANONE	<2.0	
2-HEXANONE	<10	_
TETRACHLOROETHENE	1.7	
TETRACHLOROETHENE CHLOROBENZENE ETHYLBENZENE BROMOFORM	<1.0	
ETHYLBENZENE	<1.0	
BROMOFORM	<3.0	



VOLATILE ORGANICS ANALYSIS DATA SUMMARY

_	CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : MW-7-99/00-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED : 08/24/99 DATE RECEIVED : 08/25/99 DATE EXTRACTED : N/A DATE ANALYZED : 09/06/99 UNITS : ug/L DILUTION FACTOR : 1
<u> </u>	COMPOUNDS	RESULTS
	(m+p)-XYLENE STYRENE O-XYLENE 1,1,2,2-TETRACHLOROETHANE 1,3-DICHLOROBENZENE 1,4-DICHLOROBENZENE P-ISOPROPYLTOLUENE 1,2-DICHLOROBENZENE	<1.0 <1.0 <1.0 <1.0 <2.0 <2.0 <2.0 <2.0

SURROGATE PERCENT RECOVERY		LIMITS
1,2-DICHLOROETHANE-D4 TOLUENE-D8	138 H 89 89	81 - 130 80 - 120 75 - 118

H = Out of limits.

MAS I.D. # 908046-14RE

: ERM-WEST

CLIENT

VOLATILE ORGANICS ANALYSIS DATA SUMMARY

PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : MW-7-99/00-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED : 08/24/99 DATE RECEIVED : 08/25/99 DATE EXTRACTED : N/A DATE ANALYZED : 09/08/99* UNITS : ug/L DILUTION FACTOR : 1	-
COMPOUNDS		7
CHLOROMETHANE	<5.0	_
VINYL CHLORIDE	<1.0	
BROMOMETHANE	<1.0	
VINYL CHLORIDE BROMOMETHANE CHLOROETHANE TRICHLOROETHOROMETHANE	<1.0	_
THE CHILD HOUR DOOR OFFE THANK	<1.0	
1 1-DICUIODOEMUDNU	<10	
ACETONE 1,1-DICHLOROETHENE METHYLENE CHLORIDE CARRON DISH FIRE	<1.0 <5.0	
CARBON DISULFIDE	<5.0	
TRANS-1 2-DICULODOEMUENE	<5.0	
1.1-DICHLOROFTHANE	<1.0	
CARBON DISULFIDE TRANS-1,2-DICHLOROETHENE 1,1-DICHLOROETHANE VINYL ACETATE 2-BUTANONE	<1.0	_
2-BUTANONE	<10	
	<10	
CHLOROFORM CIS-1,2-DICHLOROETHENE 1,1,1-TRICHLOROETHANE 1,2-DICHLOROETHANE	<1.0	
1,1,1-TRICHLOROETHANE	<1.0	_
1,2-DICHLOROETHANE	<1.0	ا
CARBON TETRACHLORIDE	<1.0	-
1,2-DICHLOROETHANE CARBON TETRACHLORIDE BENZENE 1,2-DICHLOROPROPANE TRICHLOROETHENE	<1.0	-
1,2-DICHLOROPROPANE	<1.0	1
TRICHLOROETHENE	2.2	
TRICHLOROETHENE BROMODICHLOROMETHANE CIS-1,3-DICHLOROPROPENE 4-METHYL-2-PENTANONE TRANS-1,3-DICHLOROPROPENE	<1.0	_
CIS-1, 3-DICHLOROPROPENE	<3.0	
4-METHYL-2-PENTANONE	<10	
TRANS-1, 3-DICHLOROPROPENE	<3.0	
1,1,2-TRICHLOROETHANE	<1.0	_
CULODODIDO CACAMINA	<1.0	
TRANS-1, 3-DICHLOROPROPENE 1, 1, 2-TRICHLOROETHANE TOLUENE CHLORODIBROMOMETHANE 2-HEXANONE	<2.0	
	<10	_
TETRACHLOROETHENE CHLOROBENZENE ETHYLBENZENE BROMOFORM	1.8	
ETHYLRENGENE	<1.0	
BROMOFORM	<1.0	
DIOPOPOM	<3.0	

RE = Reanalyzed.

^{* =} Reanalyzed outside of the recommended holding time.

VOLATILE ORGANICS ANALYSIS DATA SUMMARY

CLIENT	:	ERM-WEST	DATE	SAMPLED	:	08/24/99
PROJECT #	:	6007.23	DATE	RECEIVED	:	08/25/99
PROJECT NA		SEATTLE ANGS	DATE	EXTRACTED	:	N/A
CLIENT I.I	D. :	MW- 7- 9 9/00-1	DATE	ANALYZED	:	09/08/99*
SAMPLE MAT	TRIX:	WATER	UNITS	3	:	ug/L
EPA METHOI	D :	8260B	יוודע	TION FACTOR	٠	1

COMPOUNDS RESULTS

	(m+p)-XYLENE	<1.0
	STYRENE	<1.0
	O-XYLENE	<1.0
	1,1,2,2-TETRACHLOROETHANE	<1.0
-	1,3-DICHLOROBENZENE	<2.0
	1,4-DICHLOROBENZENE	<2.0
	P-ISOPROPYLTOLUENE	<2.0
_	1,2-DICHLOROBENZENE	<2.0
	•	-2.0

SURROGATE PERCENT RECOVERY		LIMITS
1,2-DICHLOROETHANE-D4 TOLUENE-D8BROMOFLUOROBENZENE	122 90 96	81 - 130 80 - 120 75 - 118

MAS I.D. # 908046-14RE

RE = Reanalyzed.
* = Reanalyzed outside of the recommended holding time.



COMPOUNDS RESULTS CHLOROMETHANE	CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : MW-5-99/00-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED : 08/24/99 DATE RECEIVED : 08/25/99 DATE EXTRACTED : N/A DATE ANALYZED : 09/06/99 UNITS : ug/L DILUTION FACTOR : 1	
CHLOROMETHANE VINYL CHLORIDE SEROMOMETHANE CHLOROFTHANE CHLOROFTHANE CHLOROFTHANE CHLOROFTUOROMETHANE CACETONE 1,1-DICHLOROETHENE 1,1-DICHLOROETHENE CARBON DISULFIDE CARBON CIO CARBON CIO CHLOROETHANE CIO CHLOROFORM CIO CHLOROFORM CIS-1,2-DICHLOROETHENE CIO CALOROFORM CIS-1,2-DICHLOROETHENE CIO CALOROFORM CIS-1,2-DICHLOROETHANE CARBON TETRACHLORIDE CARBON TETRACHLORIDE CARBON TETRACHLORIDE CARBON TETRACHLORIDE CARBON TETRACHLOROETHANE CIO CRICLLOROETHANE CIO CRICLLOROETHANE CIO CRICLLOROETHANE CIO CRICLLOROETHANE CIO CRICLLOROETHANE CIO CRICLLOROETHANE CIS-1,3-DICHLOROPROPENE CIO CRICLLOROETHANE CIO CRICLLOROETHANE CIS-1,3-DICHLOROPROPENE CIS-1,3-DICHLOROPROPENE CIS-1,3-DICHLOROFROPENE CIS-1,3-DICHLOROFROPENE CIS-1,3-DICHLOROFROPENE CIS-1,3-DICHLOROFROPENE CIS-1,3-DICHLOROETHANE CIS-1,0 CREATACHLOROETHANE CIS-1,0 CREATACHLOROETHANE CIS-1,0 CREATACHLOROETHANE CIS-1,0 CREATACHLOROETHANE CIS-1,0 CREATACHLOROETHANE CIS-1,0 CIS-1,0 CREATACHLOROETHANE CIS-1,0 CIS-1,0 CREATACHLOROETHANE CIS-1,0 CIS-1,0 CREATACHLOROETHANE CIS-1,0 CIS-1,0 CREATACHLOROETHANE CIS-1	COMPOUNDS	RESULTS	
~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~	CHLOROMETHANE VINYL CHLORIDE BROMOMETHANE CHLOROETHANE TRICHLOROFLUOROMETHANE ACETONE 1,1-DICHLOROETHENE METHYLENE CHLORIDE CARBON DISULFIDE TRANS-1,2-DICHLOROETHENE 1,1-DICHLOROETHANE VINYL ACETATE 2-BUTANONE CHLOROFORM CIS-1,2-DICHLOROETHANE 1,2-DICHLOROETHANE 1,2-DICHLOROETHANE 1,2-DICHLOROETHANE CARBON TETRACHLORIDE BENZENE 1,2-DICHLOROPROPANE TRICHLOROETHANE CIS-1,3-DICHLOROPROPENE 4-METHYL-2-PENTANONE TRANS-1,3-DICHLOROPROPENE 1,1,2-TRICHLOROETHANE CIS-1,2-TRICHLOROETHANE CIS-1,3-DICHLOROPROPENE 4-METHYL-2-PENTANONE TRANS-1,3-DICHLOROPROPENE 1,1,2-TRICHLOROETHANE CHLORODIBROMOMETHANE CHLORODIBROMOMETHANE CHLOROBENZENE ETHYLBENZENE ETHYLBENZENE	<5.0 <1.0 <1.0 <1.0 <1.0 <5.0 <5.0 <5.0 <1.0 <1.0 <1.0 <1.0 <1.0 <1.0 <1.0 <1	



# VOLATILE ORGANICS ANALYSIS DATA SUMMARY

_	CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : MW-5-99/00-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED DATE RECEIVED DATE EXTRACTE DATE ANALYZED UNITS DILUTION FACT	: 08/25/99 D : N/A : 09/06/99 : ug/L
<u> </u>	COMPOUNDS	RESULTS	
_	(m+p)-XYLENE STYRENE O-XYLENE 1,1,2,2-TETRACHLOROETHANE 1,3-DICHLOROBENZENE 1,4-DICHLOROBENZENE P-ISOPROPYLTOLUENE 1,2-DICHLOROBENZENE	<1.0 <1.0 <1.0 <1.0 <2.0 <2.0 <2.0 <2.0	
	SURROGATE PERCENT RECOVERY		LIMITS
	1,2-DICHLOROETHANE-D4 TOLUENE-D8	136 H 89 88	81 - 130 80 - 120 75 - 118

H = Out of limits.



CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : BS-004PZ-99/00-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B		
COMPOUNDS	RESULTS	
CHLOROMETHANE VINYL CHLORIDE BROMOMETHANE CHLOROETHANE TRICHLOROFLUOROMETHANE ACETONE 1,1-DICHLOROETHENE METHYLENE CHLORIDE CARBON DISULFIDE	<5.0 <1.0 <1.0 <1.0 <1.0 <1.0 <5.0 <5.0 <5.0 <1.0 <1.0 <1.0 <1.0 <1.0 <1.0 <1.0 <1	
	<3.0	- 1



_	CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : BS-004PZ-99/00-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED DATE RECEIVED DATE EXTRACTE DATE ANALYZED UNITS DILUTION FACT	: 08/25/99 D : N/A : 09/08/99 : ug/L
1	COMPOUNDS	RESULTS	
	(m+p)-XYLENE STYRENE O-XYLENE 1,1,2,2-TETRACHLOROETHANE 1,3-DICHLOROBENZENE 1,4-DICHLOROBENZENE P-ISOPROPYLTOLUENE 1,2-DICHLOROBENZENE	<1.0 <1.0 <1.0 <1.0 <2.0 <2.0 <2.0 <2.0	
	SURROGATE PERCENT RECOVERY		LIMITS
_	1,2-DICHLOROETHANE-D4 TOLUENE-D8 BROMOFLUOROBENZENE	123 94 96	81 - 130 80 - 120 75 - 118



EDA MERICO	DATE SAMPLED : 08/25/99 DATE RECEIVED : 08/25/99 DATE EXTRACTED : N/A DATE ANALYZED : 09/08/99 UNITS : ug/L DILUTION FACTOR : 1	
COMPOUNDS	RESULTS	-
CHLOROMETHANE VINYL CHLORIDE BROMOMETHANE CHLOROETHANE TRICHLOROFLUOROMETHANE ACETONE 1,1-DICHLOROETHENE METHYLENE CHLORIDE CARBON DISULFIDE TRANS-1,2-DICHLOROETHENE 1,1-DICHLOROETHANE VINYL ACETATE 2-BUTANOME	<5.0 <1.0 <1.0 <1.0 <1.0 <1.0 <5.0 <5.0 <5.0 <1.0 <1.0 <1.0 <1.0 <1.0 <1.0 <1.0 <1	
BROMOFORM	<1.0 <3.0	



-	CLIENT : ERM-WEST PROJECT # : 6007.23 PROJECT NAME : SEATTLE ANGS CLIENT I.D. : MW-10-99/00-1 SAMPLE MATRIX : WATER EPA METHOD : 8260B	DATE SAMPLED DATE RECEIVED DATE EXTRACTE DATE ANALYZED UNITS DILUTION FACT	: 08/25/99 D : N/A : 09/08/99 : ug/L
<u> </u>	COMPOUNDS	RESULTS	
	(m+p)-XYLENE STYRENE O-XYLENE 1,1,2,2-TETRACHLOROETHANE 1,3-DICHLOROBENZENE 1,4-DICHLOROBENZENE P-ISOPROPYLTOLUENE 1,2-DICHLOROBENZENE	<1.0 <1.0 <1.0 <1.0 <2.0 <2.0 <2.0 <2.0	
_	SURROGATE PERCENT RECOVERY		LIMITS
	1,2-DICHLOROETHANE-D4 TOLUENE-D8	114 90 93	81 - 130 80 - 120 75 - 118



CLIENT PROJECT # PROJECT NAME CLIENT I.D. SAMPLE MATRIX EPA METHOD	: MW-1-99/00-1	DATE SAMPLED : 08/25/99 DATE RECEIVED : 08/25/99 DATE EXTRACTED : N/A DATE ANALYZED : 09/08/99 UNITS : ug/L DILUTION FACTOR : 1
--------------------------------------------------------------------	----------------	---------------------------------------------------------------------------------------------------------------------------------

<pre>RESULTS  &lt;5.0 &lt;1.0 &lt;1.0 &lt;1.0 &lt;1.0 &lt;1.0 &lt;1.0 &lt;1.0 &lt;1</pre>
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<1.0 <1.0 <1.0 <1.0 <1.0 <5.0 <5.0 <1.0 <1.0 <1.0 <1.0
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4



### VOLATILE ORGANICS ANALYSIS DATA SUMMARY

CLIENT : ERM-WEST DATE SAMPLED : 08/25/99 PROJECT #
PROJECT NAME : 6007.23 DATE RECEIVED : 08/25/99 : SEATTLE ANGS DATE EXTRACTED : N/A CLIENT I.D. : MW-1-99/00-1DATE ANALYZED : 09/08/99 SAMPLE MATRIX : WATER UNITS : ug/L EPA METHOD : 8260B

	200B	DILUTION FACTOR: 1
<u> </u>	COMPOUNDS	RESULTS
	(m+p)-XYLENE STYRENE O-XYLENE 1,1,2,2-TETRACHLOROETHANE 1,3-DICHLOROBENZENE 1,4-DICHLOROBENZENE P-ISOPROPYLTOLUENE 1,2-DICHLOROBENZENE	<1.0 <1.0 <1.0 <1.0 <2.0 <2.0 <2.0 <2.0
	SURROGATE PERCENT RECOVERY	LIMITS
	1,2-DICHLOROETHANE-D4 TOLUENE-D8 BROMOFLUOROBENZENE	115 93 96 81 - 130 80 - 120 96 75 - 118

## APPENDIX C

# QUALITY CONTROL DATA VALIDATION REPORT-AUGUST 1999 GROUNDWATER DATA

### **APPENDIX C**

# QUALITY CONTROL DATA VALIDATION REPORT - AUGUST 1999 GROUNDWATER DATA

Analytical data are the basis for evaluating the environmental conditions at the Seattle Air National Guard Station (Seattle ANGS) in Seattle, Washington. A primary objective of environmental sampling conducted at the site is to obtain accurate data that reflect actual conditions.

This report addresses groundwater analytical data collected in August 1999 at the Seattle ANGS as part of a quarterly monitoring program. Eighteen groundwater samples were analyzed for volatile organic analytes using United States Environmental Protection Agency (USEPA) Method 8260. To ensure that data quality was acceptable for decision-making purposes, analytical data for this project were validated. This process identifies limitations on the use of the data, or data that should not be used for decision-making purposes. The quality of the data was assessed and any necessary qualifiers were applied following the USEPA Contract Laboratory Program National Functional Guidelines for Organic Data Review (February 1994).

Environmental Resources Management (ERM) validated data for compliance with the following quality assurance/quality control (QA/QC) project- and/or method-prescribed criteria:

- Holding Time and Preservation: The holding time is the period of time between collection of the sample and preparation/analysis of the sample. Analyses performed for this project have method-prescribed holding times. Preservation refers to the temperature at which the samples are received at the laboratory, as well as any pH anomalies noted by the laboratory for acid-preserved samples.
- Calibration: The analysis of target analytes at a range of concentrations to develop a graphical plot of instrument response against the different analyte concentrations. An initial calibration

curve establishes the graphical plot, and the continuing calibration verification monitors daily instrument linearity against the initial calibration.

- Blank Samples: The preparation and analysis of samples from reagent (contaminant-free) water. Blank samples for this investigation included method, trip, rinsate, and field blanks. Detections in a blank sample indicate laboratory, handling, or field contamination.
- Internal Standards: The addition of compounds similar to target analytes of interest that are added to sample aliquots for organic analysis. The internal standards are used to quantitatively and qualitatively evaluate retention time and instrument response for each analytical run.
- Spike Samples: The preparation and analysis of an environmental sample or a sample of reagent water spiked with a subset of target analytes at known concentrations. The results of the spike analysis measure laboratory accuracy in the reagent sample, and results from the environmental sample spike measure potential interferences from the sample matrix.
- Surrogate Spikes: The addition of compounds similar to target analytes of interest that are added to sample aliquots for organic analysis. Surrogate spikes measure possible interferences from the sample matrix for the analysis of target analytes.
- Duplicate Samples: The preparation and analysis of an additional aliquot of the sample. The results from duplicate analysis measure potential heterogeneity of contaminant concentrations in the samples.

The following data qualifiers were used as appropriate during this validation process:

- U: The analyte was analyzed for, but was not detected above the reported quantitation limit.
- J: The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- UJ: The analyte was not detected above the reported sample quantitation limit; however, the reported quantitation limit was approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.

R: The sample results were rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte could not be verified.

None of the August 1999 groundwater data were rejected based on the data validation. All of the data, including qualified data, are acceptable and can be used for decision-making purposes.

The following discussion addresses the results of the data validation for each of the QA/QC components listed above.

### **Holding Time and Preservation**

The USEPA has established maximum recommended holding times for the analyses performed on the August 1999 groundwater samples. The USEPA has also defined the acceptable temperature range at which samples must be stored to maintain sample preservation. Holding times and sample temperatures extending beyond the maximum can negatively affect sample integrity (e.g., loss of volatile compounds, biodegradation), and impacted samples are qualified depending on the severity of the exceedence and the analytes of concern. The maximum recommended holding time for USEPA Method 8260 is 14 days for acid-preserved samples and 7 days for unpreserved samples. The recommended temperature range for sample storage is 2 to 6 degrees Celsius.

Each of the sample analysis results was reviewed for compliance with the method-prescribed preparation and analysis holding times. None of the initial analyses were performed outside of the prescribed holding time. Accordingly, initial sample results were not compromised by an excessive period between sample collection and analysis. Three of the samples required reanalysis due to elevated surrogate recoveries. The reanalysis was performed one day past the recommended holding time of 14 days. The reanalysis results were qualified "J," estimated concentration, and "UJ," estimated reporting limit, for detected and non-detected (ND) analytes. The affected data are listed in Table C-1.

The temperature of the samples upon receipt at the laboratory was also reviewed for compliance with method requirements. The samples were received within the method-prescribed temperature range. None of the samples were compromised by out-of-range temperature preservation.

### **FINAL**

### TABLE C-1

Data Qualified Based on Exceeded Holding Time 143rd CCSQ, Seattle ANGS, Seattle, Washington

Data	Sample ID	Analytical	Number	ERM
Package	<del></del>	Parameter	of days exceeded	Qualifier*
908046	MW-8-99/00-1 RE	VOCs	1	]/ប្យ
	MW-6-99/00-1 RE	VOCs	1	j/Uj
	MW-7-99/00-1 RE	VOCs	1	J/UJ

#### Notes:

RE = Reanalysis of sample

VOC = Volatile organic compound

J = Reported values for detected analytes are estimated concentrations.

UJ = Reported detection limits for analytes are estimated concentrations.

^{*}Data qualifiers apply to listed samples.

### **Calibration Results**

Before an analytical instrument is used for sample analysis, the instrument should be calibrated to within USEPA method specifications. The purpose of this calibration is to ensure that the instrument is appropriately responsive to measurable chemical concentrations. If an instrument is not properly calibrated, it may not be capable of producing acceptable quantitative, qualitative, and reproducible data. For example, positive detections of a given analyte could contain an undetermined degree of inaccuracy if the instrument is out of calibration, although the results may still be considered valid. In the case of ND analytes, the associated reporting limit would be similarly affected; however, such results would still be considered NDs.

Two types of calibration data were reviewed: initial calibration verification (ICV) and continuing calibration verification (CCV). The ICV consisted of standards that were analyzed at five or more concentrations. These concentrations ranged from the reporting limit to the upper linear range of the instrument. Average response factors from the ICV were used to calculate sample results. The laboratory evaluated the ICVs using relative standard deviation (RSD). The reported RSDs were compared to the method-prescribed acceptance criteria during the data validation. None of the ICV RSDs exceeded the acceptable method-prescribed criteria. Accordingly, none of the data were qualified based on the initial calibration results.

The CCV is analyzed either daily or every 12 hours to ensure the instrument response is still within method performance criteria for linearity. The CCV consisted of analyzing a standard at one concentration; the concentration of this standard was generally in the midrange of the ICV standard concentrations. The laboratory calculated the percent difference (%D) between the CCV and the ICV. The %Ds were compared to the method-prescribed acceptance criteria during the data validation. Table C-2 lists the CCV %Ds that were above the method-prescribed criteria and the samples associated with each CCV, along with the applied data qualifiers. None of the analytes associated with an elevated %D were detected in the project samples.

### **Blank Samples**

Blank samples are prepared in the laboratory or in the field and are carried through the analytical process. The purpose of a blank sample is

**FINAL** 

TABLE C-2

Data Qualified Based on Calibration Results
143rd CCSQ, Seattle ANGS, Seattle, Washington

Analytical	Calibration	Associated	Instrument	Date	Target Compound	RSD	ERM
Parameter	(ICV/CCV)	Samples	ID			or %D	Qualifier
VOCs	CCV	TB082499-1	HP4	9/2/99 am	Trichlorofluoromethane	31.1	UI
		MW-8-99/00-1FA			Vinyl Acetate	50.7	บั
		MW-8-99/00-1FT			Carbon tetrachloride	49.6	ญ่
		BS-006PZ-99/00-1			Chlorodibromomethane	39.4	ហ៍
		MW-2-99/00-1			Bromoform	48.6	υj
	CCV	BS-005PA-99/00-1	HP4	9/2/99 pm	Trichlorofluoromethane	31	ບງ
		MW-9-99/00-1		•	Carbon disulfide	33.5	บ
		MW-2-99/00-1			Vinyl Acetate	<b>45</b> .1	บ
					2-Butanone	39	ÚĴ
					Carbon tetrachloride	50.8	ບັງ
	CCV	MW-4-99/00-1	HP4	09/05/1999	Chloromethane	31	បា
		MW-4-99/00-1R			Trichlorofluoromethane	32	ຫຼ
		MW-8-99/00-1			Vinyl Acetate	47.6	UJ
		MW-8-99/00-1D			Carbon tetrachloride	59.8	ຫຼ
		MW-6-99/00-1			Chlorodibromomethane	58	UJ
		MW-7-99/00-1			Bromoform	<i>7</i> 7.6	UJ
		MW-5-99/00-1					-
	CCV	MW-8-99/00-1D R	HP4	09/08/1999	Trichlorofluoromethane	39.7	UJ
		MW-6-99/00-1 RE			Acetone	31.2	ÚJ
		MW-7-99/00-1 RE			Vinyl acetate	50.1	UJ
		BS-004PZ-99/00-1			1,2-Dichloroethane	60.3	υj
		MW-10-99/00-1			Bromodichloromethane	33.9	UJ
		MW-1-99/00-1			Chlorodibromomethane	53.8	υj
					Bromoform	73.8	UJ

#### Notes:

UJ = Reported detection limits for the listed compounds and samples are estimated concentrations.

TB = Trip blank sample

FA = Field blank sample prepared with ASTM Type II water

FT = Field blank sample prepared with tap water

D = Duplicate sample

RE = Reanalysis of sample

RSD = Relative standard deviation

[%]D = Percent difference

VOC = Volatile organic compound

ICV= Initial calibration verification

CCV= Continuing calibration verification

^{*} Data qualifiers apply to all related samples

to test for contamination resulting from laboratory, shipping, or other sample-handling activities. Blank samples are analyzed and evaluated for detections of target analytes. If target analytes are detected in a blank sample, these detections indicate that some element of the sample collection or analysis process has introduced contaminants not present in the original environmental sample aliquot. If target analytes are detected in a blank sample, then all associated data must be evaluated to determine whether:

- Those data have been similarly impacted; or
- The blank detections are an isolated occurrence not representative of other data.

The four types of blank samples analyzed and reported with the groundwater samples collected in August 1999 were method, trip, rinsate, and field blank samples. Preparation, handling, and analysis of these blank samples are summarized below.

- Method blank samples monitor for potential laboratory contamination of samples. Method blank samples were prepared in the laboratory by taking an aliquot of reagent water through all preparation and analysis steps. A method blank was prepared and analyzed with each batch of environmental samples.
- 2. Trip blank samples monitor for potential contamination of samples during collection and transportation to the laboratory. Trip blank samples were prepared by filling a volatile organics analysis (VOA) vial with an aliquot of reagent water and sealing it with a Teflon-lined-septum lid. The trip blank sample travels with the filled aqueous sample containers to the laboratory.
- 3. Rinsate blank samples monitor for potential contamination of project samples from inadequate decontamination of sample collection equipment. Rinsate blank samples were prepared in the field by pouring American Society for Testing and Materials (ASTM) Type II reagent-grade water over the decontaminated sample collection equipment. The water was collected in clean sample containers supplied by the laboratory. Rinsate blank samples were labeled with an "R" identifier at the end of the sample ID.
- 4. Field blank samples monitor for potential contamination of project samples from ambient conditions at the sample collection site. Field blank samples were prepared at sample collection locations by slowly

pouring tap water or ASTM Type II water into clean sample containers supplied by the laboratory. Field blank samples prepared with tap water were labeled with an "FT" identifier at the end of the sample ID. The identifier "FA" was used to designate field blank samples prepared with ASTM Type II water.

No target analytes were detected in the method, trip, or rinsate blank samples. The common drinking water contaminants chloroform and bromodichloromethane were reported the tap water field blank sample. Three of the associated sample results required qualification based on the tap water field blank results. The sample results were qualified as ND following the USEPA National Functional Guidelines 5-times rule. Table C-3 lists the field blank and associated sample-detected concentrations, along with the applied data qualifiers.

### Spike Samples

A spike sample is a QC sample that is prepared and analyzed by the laboratory. The laboratory prepares, analyzes, and reports spike sample results to demonstrate their ability to properly analyze, detect, and quantify target analytes. A spike sample result is typically reported as the amount of analyte detected divided by the known amount spiked into the sample, and is commonly referred to as percent recovery. The percent recovery is then compared to an established limit range. ² The two types of spike samples analyzed with the project samples were matrix and blank spikes.

 Matrix spike (MS) samples consist of an aliquot of an environmental sample that is spiked with known concentrations of a subset of target analytes. A matrix spike duplicate (MSD) sample is a second (duplicate) spike sample prepared and analyzed with the MS sample. MS samples are used to monitor potential interference from the sample matrix for target analytes. A low MS recovery may indicate lowbiased sample results; a high MS recovery may indicate high-biased sample results.

¹ The USEPA CLP National Functional Guidelines state "any compound detected in the sample (other than common laboratory contaminants) that was also detected in any associated blank is qualified if the sample concentration is less than five times (5x) the blank concentration."

² In most cases, the prescribed analytical method will specify protocol to develop appropriate limit ranges. In some cases, however, limit ranges are established by the laboratory in the method procedures.

TABLE C-3
Blank Detections and Associated Qualifed Sample Results
143rd CCSQ, Seattle ANGS, Seattle, Washington

Data	Blank	Sample ID	Target	Reported	ERM
908046	Sample ID		Compound	Concentration (µg/l)	Qualifier*
908046	MW-8-8-99/00-1FT		Chloroform	45	
			Bromodichloromethane	1.6	
		BS-004PZ-99/00-1	Chloroform	65	U
			<b>Bromodichloromethane</b>	2.3	Ü
		MW-10-99/00-1	Chloroform	63	U
			Bromodichloromethane	2.3	Ü
		MW-1-99/00-1	Chloroform	67	Ü
			Bromodichloromethane	2.3	Ū

#### Notes:

FT = Field blank sample prepared with tap water

U = Reported detections should be considered non-detected at the reported concentration.

^{*}Data qualifiers apply to listed samples.

2. Blank spike samples, commonly referred to as laboratory control samples (LCS), consist of an aliquot of reagent water that is spiked with known concentrations of a subset of target analytes. The LCS sample is used to monitor laboratory accuracy without the bias of a sample matrix. LCS recoveries outside of acceptable limits may indicate poor laboratory accuracy.

All of the MS and LCS recoveries were within acceptable limits. The acceptable MS and LCS recoveries indicate minimal matrix interference and acceptable laboratory accuracy for the August 1999 groundwater data.

### **Internal Standard Responses**

Under USEPA methods, a given analyte list for organic compounds is segregated by chemical properties and retention time into one or more subsets. A USEPA-defined internal standard with comparable chemical properties and retention times is assigned to each subset of analytes. The laboratory adds a known concentration of an internal standard to each sample, including laboratory QC samples (e.g., calibration standards, MS, method blank samples), prior to analysis. The instrument internal standard response for each sample is compared to the internal standard response in the daily CCV. The sample internal standard area count must be within the range of 0.5 to 2 times the CCV area count, and the retention time must be within ±30 seconds of the CCV retention time. If the area count and/or retention time measured for the sample is outside the acceptable range, quantitation results for the associated analyte subset may be biased. Interferences from the sample matrix are typically responsible for internal standard responses that are consistently outside acceptable ranges; most matrix interferences cause a consistently high or low bias.

Internal standards were added to each of the project samples analyzed for VOCs. The internal standard responses were within acceptable limits, indicating minimal matrix interference and acceptable sample quantitation.

### Surrogate Spikes

A surrogate spike is similar to an internal standard; it is chemically similar to the target analytes and is only used in organic analyses. The difference between surrogate spikes and internal standards is that surrogate spikes

are used only to assess possible interferences from the sample matrix, whereas internal standards are used to quantitate target analytes while accounting for any interferences from the sample matrix. Surrogate spike results are typically reported in terms of percent recovery, which is calculated by dividing the amount of surrogate detected in the sample by the known amount of surrogate added to the sample.

For the August 1999 groundwater data, surrogate recoveries were compared to the laboratory-generated limits of acceptance. The surrogate recoveries were mostly within acceptable limits, indicating that sample results were subject to minimal interferences from the sample matrix.

Five of the August 1999 samples had one slightly elevated surrogate recovery. Positive detections in these five samples were qualified "J," estimated concentration. The laboratory reanalyzed three of the five samples with positive detections to confirm the elevated surrogate recovery. The surrogate recoveries in the reanalysis were within acceptable limits; however the reanalysis was performed one day past the recommended holding time. Samples with out-of-limit surrogate recoveries and the applied data qualifiers are listed in Table C-4.

### **Duplicate Samples**

A duplicate sample is a second aliquot of a sample that is treated the same as the primary sample. A duplicate sample analysis is performed to measure the precision of the method and possible heterogeneity of analyte concentrations in the sample matrix. Duplicate field samples are collected to measure matrix heterogeneity.

Laboratory duplicate analyses for the project samples consisted of matrix spike duplicate analyses. The laboratory calculated the relative percent difference (RPD) between the MS and MSD spike concentrations. The calculated RPDs were compared to method-prescribed or laboratory-generated acceptable limits. A field duplicate sample also was collected and submitted for analysis, and an RPD was calculated for detected analytes.

The duplicate sample RPDs were within acceptable limits, indicating acceptable analytical precision and minimal matrix heterogeneity.

### **FINAL**

TABLE C-4
Surrogate Recovery Results Outside of Acceptable Limits
143rd CCSQ, Seattle ANGS, Seattle, Washington

Data	Sample ID	Analytical	Surrogate	% Recovery	Control	ERM
Package 908046		Parameter			Limit	Qualifier
908046	MW-4-99/00-1R	VOCs	1,2-Dichloroethane-d4	133	81-130	
	MW-8-99/00-1D	VOCs	1,2-Dichloroethane-d4	136	81-130	1
	MW-6-99/00-1	VOCs	1,2-Dichloroethane-d4	137	81-130	ĺ
	MW-7-99/00-1	VOCs	1,2-Dichloroethane-d4	138	81-130	í
	MW-5-99/00-1	VOCs	1,2-Dichloroethane-d4	136	81-130	•
	MW-8-99/00-1MS	VOCs	1,2-Dichloroethane-d4	134	81-130	
	MW-8-99/00-1MSD	<b>VOCs</b>	1,2-Dichloroethane-d4	134	81-130	

#### Notes:

R = Rinsate blank

D = Duplicate sample

MS= Matrix spike

MSD= Matrix spike duplicate

VOC = Volatile organic compound

J = Reported values for detected analytes associated with the listed surrogate are estimated concentrations.

^{*}Data qualifiers apply to target analyte detections in listed samples.

### Overall Assessment

None of the Seattle ANGS analytical data for groundwater samples collected in August 1999 were rejected. The data set is 100 percent complete and meets the project goal for completeness. Unqualified data can be used for decision-making purposes. Qualified data also can be used for decision-making purposes; however, the limitations identified by the qualifiers should be considered when using the data. The quality of the August 1999 groundwater analytical data is acceptable for the preparation of technically defensible documents.

# APPENDIX D

# CHAIN-OF-CUSTODY FORMS

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